सुगन्ध एवं सुरस विकास केन्द्र, कन्नौज
FRAGRANCE & FLAVOUR DEVELOPMENT CENTRE, KANNAUJ

भारत सरकार की एक स्वायत्तशासी संस्था
सूक्ष्म, लघु एवं मध्यम उद्यम मंत्रालय
A Govt. of India Autonomous Body
Ministry of Micro, Small & Medium Enterprises

An ISO 9001:2015, ISO 14001:2015 & ISO 29990:2010 Certified Organisation

T-0155

Quality Assessment Laboratory

TEST-REPORT

Serial No.63	Ref. No. FFDC/KNJ/QAL/2022-23/	Date: 10.05.2022

A. SAMPLE DETAILS

Customer's Name & Address	**KRAMA AGRO,** Kukde village,Boisar E Dist.Palghar,401501 Maharashtra state.
Customer's Code	
Item Name	Rose Absolute
Specification	QAL-SOPM 201
Job Order No.	58, XXXI-58
Date of receipt	05.05.2022
Date of Analysis	10.05.2022
Cash receipt No. & Date	12/1188 dt- 12/05/22

S. No.	Name Of Testing	Test Value	Test Method & Remark
1	Colour, Odour & Appearance	Reddish Brown & Petalic rose	
1	Optical rotation	Undermine	IS 326 (Part 4): 2005 (RA 2015)
2	Specific Gravity at 27 °C	0.9407	IS 326 (Part 3): 2006 (RA 2012)
3	Refractive Index at 27 °C	1.4866	IS 326 (Part 5): 2006 (RA 2012)
4	Solubility in 90% EtOH at 27°C	Soluble upto 8 Vol.	IS 326 (Part 6): 2005 (RA 2015)
5	**GLC**	Chromatogram enclosed	IS 326 (Part 19): 1998 (RA 2014)
6	GC-MS	Chromatograph Enclosed.	IS: 236 (Part 19) 1998
	The identification of compounds in said report is based upon library search software of GC-MS using NIST (USA)/Adam/ Flavour library. The Institute holds no responsibility legal or otherwise towards authenticity of identification accomplished by GC-MS.		

GLC

Disclaimer:

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2-This report shall not be reproduced except in full, without the written approval of the Centre.
3-The results are valid at the time of test and under stated conditions.
4- The results are not valid for legal implications in the court of law.
5-Any complaint may be communicated to the customer service cell, FFDC Kannauj. If we don't receive any customer feedback form it will be assumed that party is satisfied with report.

10/5/2022

Head
Customer Service Cell

10/5/2022

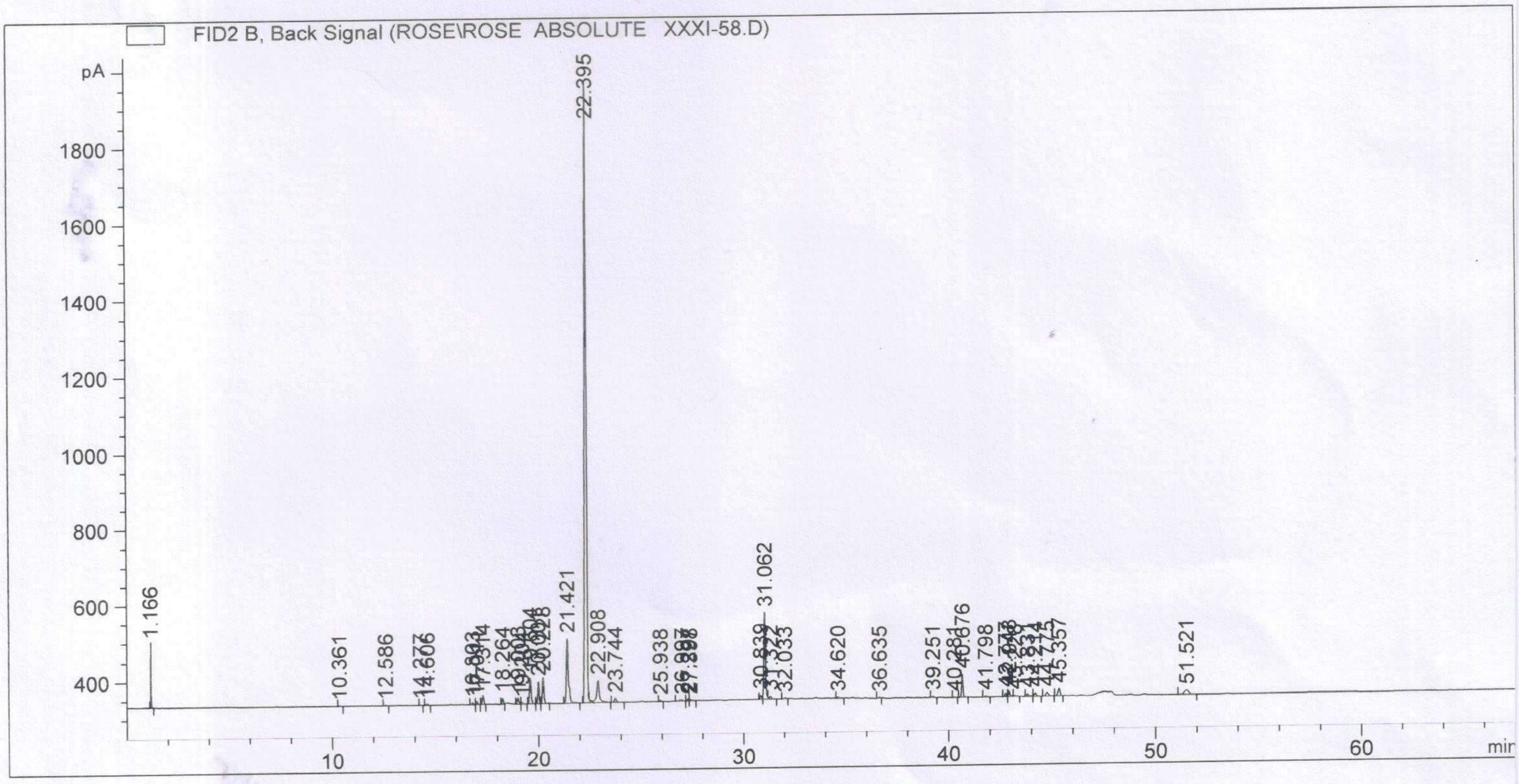
Auth. Signatory
Quality Manager/Dy. Quality Manager

Data File C:\Users\Public\Documents\ChemStation\1\Data\ROSE\ROSE ABSOLUTE XXXI-58.D
Sample Name: ROSE ABSOLUTE XXXI-58

==========

Acq. Operator : SYSTEM
Sample Operator : SYSTEM
Acq. Instrument : GC
Location : 2 (B)
Injection Date : 06-05-2022 13:16:59
Inj : 1
Inj Volume : Manually

Acq. Method : C:\Users\Public\Documents\ChemStation\1\Methods\GULAB.M
Last changed : 06-05-2022 13:12:05 by SYSTEM
(modified after loading)
Analysis Method : C:\Users\Public\Documents\ChemStation\1\Methods\GULAB.M
Last changed : 06-05-2022 14:32:38 by SYSTEM
(modified after loading)
Sample Info : ROSE ABSOLUTE XXXI-58



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Area Percent Report

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Sorted By : Signal
Multiplier : 1.0000
Dilution : 1.0000
Use Multiplier & Dilution Factor with ISTDs

Signal 1: FID2 B, Back Signal

Peak #	RetTime [min]	Type	Width [min]	Area [pA*s]	Height [pA]	Area %
1	1.166	BB	0.0159	168.61328	164.86189	1.05798
2	10.361	BB	0.0717	9.42192	1.72477	0.05912
3	12.586	BB	0.0787	12.32902	2.09422	0.07736
4	14.277	BB	0.0682	7.12176	1.50464	0.04469
5	14.606	BB	0.0813	23.19113	3.49830	0.14551

Peak #	RetTime [min]	Type	Width [min]	Area [pA*s]	Height [pA]	Area %
6	16.803	BV	0.0741	27.11535	5.44148	0.17014
7	17.011	VB	0.0715	34.81744	7.74304	0.21847
8	17.314	BB	0.0783	131.98308	25.13332	0.82814
9	18.264	BB	0.0680	69.74111	15.96949	0.43760
10	19.003	BV	0.0803	91.82078	17.20080	0.57614
11	19.204	VB	0.0834	20.72816	3.58660	0.13006
12	19.604	BB	0.0871	396.60297	67.95433	2.48852
13	20.001	BB	0.0675	251.86378	58.12908	1.58034
14	20.228	BB	0.0838	369.40726	68.57866	2.31788
15	21.421	BB	0.0998	1178.25037	167.95325	7.39304
16	22.395	BV R	0.0854	9871.54688	1613.93628	61.93993
17	22.908	VB E	0.1396	600.12537	59.01072	3.76554
18	23.744	BB	0.0965	71.92160	11.10921	0.45128
19	25.938	BB	0.0800	21.04026	4.09066	0.13202
20	26.997	BB	0.1568	45.27824	3.91922	0.28410
21	27.297	BV	0.0750	28.27919	5.78845	0.17744
22	27.398	VB	0.1340	38.42818	3.79959	0.24112
23	30.839	BV E	0.0803	60.42536	12.08307	0.37914
24	31.062	VV R	0.0885	1273.88330	227.00711	7.99310
25	31.322	VB E	0.0989	52.72549	7.51136	0.33083
26	32.033	BB	0.1015	18.29449	2.52438	0.11479
27	34.620	BB	0.0995	22.44369	2.89249	0.14083
28	36.635	BB	0.0811	11.30251	1.93645	0.07092
29	39.251	BB	0.0950	25.87331	4.19388	0.16234
30	40.281	BB	0.0802	7.15289	1.38730	0.04488
31	40.676	BB	0.0732	291.29102	59.41877	1.82773
32	41.798	BB	0.1077	28.41405	3.53890	0.17829
33	42.777	BV	0.0738	48.11269	9.87767	0.30189
34	42.943	VV R	0.0808	83.63708	15.28674	0.52479
35	43.228	VB E	0.0993	21.62368	2.92443	0.13568
36	43.831	BV	0.1189	60.11331	7.09194	0.37719
37	44.214	VB	0.1103	68.60907	9.06702	0.43049
38	44.775	BB	0.1372	87.72990	9.64172	0.55047
39	45.357	BB	0.1152	146.51181	18.95784	0.91930
40	51.521	BB	0.1974	159.51898	10.31390	1.00092
Totals :				1.59373e4	2718.68297	

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*** End of Report ***

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Specification	QAL-SOPM 201
Job Order No.	58, XXXI-58
Date of receipt	05.05.2022
Date of Analysis	10.05.2022
Cash receipt No. & Date	12/1188 dt- 12/05/22

S. No.	Name Of Testing	Test Value	Test Method & Remark
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GCMS

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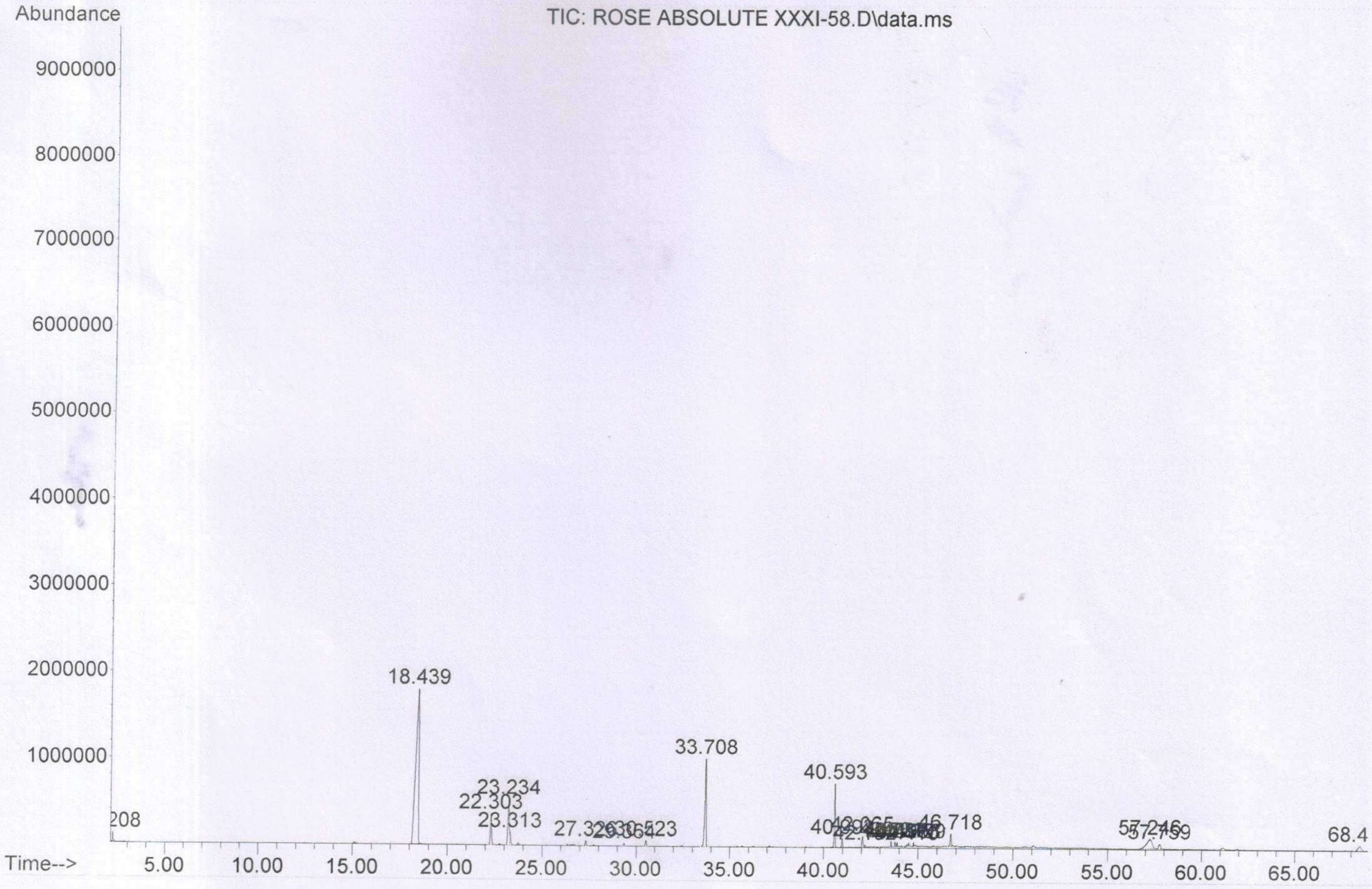
10/05/22

Head
Customer Service Cell

10/5/2022

Auth. Signatory
Quality Manager/Dy. Quality Manager

File :D:\data\New Folder\QAL feb.2016\ROSE ABSOLUTE XXXI-58.D
Operator :
Acquired : 9 May 2022 15:32 using AcqMethod PERFUME---.M
Instrument : GCMS
Sample Name: ROSE ABSOLUTE XXXI-58
Misc Info :
Vial Number: 1



Library Search Report

Data Path : D:\data\New Folder\QAL feb.2016\
Data File : ROSE ABSOLUTE XXXI-58.D
Acq On : 9 May 2022 15:32
Operator :
Sample : ROSE ABSOLUTE XXXI-58
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\DEMO1.L\Adam.L Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: ChemStation Integrator - autoint1.e

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
1	2.210	0.54	C:\Database\DEMO1.L\Adam.L			
		7.60	OCTANOL<2->	37	000123-96-6	2
		3.03	BUTANEDIOL<2,3->	337	000513-85-9	1
2	18.438	51.61	C:\Database\DEMO1.L\Adam.L			
		11.75	PHENYL ETHYL ALCOHOL (=2-PHENYLETHYL ALCOHOL;	1406	000060-12-8	91
		20.64	ISOAMYL BENZYL ETHER (=ISOPENTYL BENZYL ETHE	878	000122-73-6	42
		10.15	TOLUALDEHYDE<META->	1554	000620-23-5	9
3	22.305	4.27	C:\Database\DEMO1.L\Adam.L			
		16.80	CITRONELLOL (=3,7-DIMETHYL-6-OCTEN-1-OL)	1407	000106-22-9	96
		17.95	GERANIOL	755	000106-24-1	53
		33.15	SESQUILAVANDULOL<Z->	1453	131485-70-6	42
4	23.235	6.43	C:\Database\DEMO1.L\Adam.L			
		16.97	NEROL	1175	000106-25-2	64
		19.32	NERYL FORMATE	1183	002142-94-1	59
		34.10	SESQUILAVANDULOL<E->	1456	120707-27-9	56
5	23.310	1.32	C:\Database\DEMO1.L\Adam.L			
		24.31	PHENYL ETHYL ISOBUTANOATE	1329	007775-39-5	72
		26.27	PHENYL ETHYL BUTANOATE<2->	1319	000103-52-6	50
		18.20	PHENYL ETHYL ACETATE<2-> (=BETA-PHENYL ETHYL	1316	000103-45-7	50
6	27.327	0.61	C:\Database\DEMO1.L\Adam.L			
		19.70	LAVANDULYL ACETATE	1188	020777-39-3	91
		22.81	NERYL ACETATE	1027	000141-12-8	72
		27.82	GERANYL PROPANOATE	768	000105-90-8	72
7	29.365	0.36	C:\Database\DEMO1.L\Adam.L			
		19.18	ETHYL CHRYSANTHEMUMATE<TRANS->	690	041641-25-2	30
		34.48	THUJOPSANONE<3-ISO->	586	025966-81-8	28
		34.95	THUJOPSANONE<3->	583	025966-79-4	23
8	30.520	0.74	C:\Database\DEMO1.L\Adam.L			
		28.15	GERMACRENE D	520	023986-74-5	98
		28.53	MUUROLA-4(14),5-DIENE<TRANS-	530	000000-00-0	93
		24.04	CUBEBENE<BETA->	486	013744-15-5	93
9	40.592	5.82	C:\Database\DEMO1.L\Adam.L			
		43.03	HEXADECANOL<N->	34	036653-82-4	90
		39.42	PENTADECANOL<N->	31	000629-76-5	78
		32.44	HEXADECENE<1->	29	000629-73-2	70

```
10  40.993  0.86 C:\Database\DEMO1.L\Adam.L
                 28.82 PENTADECANE                          100 000629-62-9 86
                 20.19 TRIDECANE                             98 000629-50-5 64
                 24.61 TETRADECANE                           99 000629-59-4 64

11  42.068  1.36 C:\Database\DEMO1.L\Adam.L
                 18.77 NONANOIC ACID                       1203 000112-05-0 16
                 31.47 DODECANOIC ACID  (=LAURIC AC         648 000143-07-7 12
                 ID)
                 6.05 VALERIC ACID<3-METHYL-> (=3-M        1578 000105-43-1 10
                 ETHYLPENTANOIC AC

12  42.191  0.25 C:\Database\DEMO1.L\Adam.L
                 32.44 HEXADECENE<1->                        29 000629-73-2 58
                 40.02 OCTADECENE<1->                      1258 000112-88-9 58
                 53.22 DOCOSENE<1->                        1260 001599-67-3 58

13  43.582  0.38 C:\Database\DEMO1.L\Adam.L
                 43.03 HEXADECANOL<N->                       34 036653-82-4 86
                 46.92 EICOSENE<1->                        1259 003452-07-1 70
                 40.02 OCTADECENE<1->                      1258 000112-88-9 70

14  43.785  0.39 C:\Database\DEMO1.L\Adam.L
                 40.02 OCTADECENE<1->                      1258 000112-88-9 90
                 46.92 EICOSENE<1->                        1259 003452-07-1 90
                 53.22 DOCOSENE<1->                        1260 001599-67-3 87

15  43.897  0.37 C:\Database\DEMO1.L\Adam.L
                 32.89 HEXADECANE                          101 000544-76-3 53
                 36.74 HEPTADECANE                         102 000629-78-7 53
                 24.61 TETRADECANE                          99 000629-59-4 52

16  44.411  0.11 C:\Database\DEMO1.L\Adam.L
                 50.33 METHYL LINOLEATE (=METHYL (9        1139 000112-63-0 53
                 Z,12Z)-OCTADECADI
                 44.92 MUSK AMBRETTE (=MUSK NATURAL        1151 000123-69-3 40
                 ; =AMBRETTOLID, =
                 28.07 CITRONELLOL ISOBUTANOATE (=C         473 000097-89-2 38
                 ITRONELLYL ISOBUT

17  44.507  0.26 C:\Database\DEMO1.L\Adam.L
                 37.80 LONGIFOLOL<ISO->                    1062 001139-17-9 38
                 19.00 MENTH-1-EN-7-AL<PARA->              1465 000000-00-0 33
                 13.10 LIMONENE OXIDE<TRANS->              1040 004959-35-7 32

18  44.769  0.29 C:\Database\DEMO1.L\Adam.L
                 31.47 DODECANOIC ACID  (=LAURIC AC         648 000143-07-7 27
                 ID)
                 46.92 EICOSENE<1->                        1259 003452-07-1 18
                 40.02 OCTADECENE<1->                      1258 000112-88-9 18

19  46.716  1.36 C:\Database\DEMO1.L\Adam.L
                 40.02 OCTADECENE<1->                      1258 000112-88-9 90
                 46.92 EICOSENE<1->                        1259 003452-07-1 90
                 49.76 OCTADECANOL<N->                       35 000112-92-5 86

20  57.247  6.18 C:\Database\DEMO1.L\Adam.L
                 28.05 VANILLONE, ACETO-                   1581 000498-02-2  2
                 24.32 VANILLIN  (=BENZALDEHYDE, 4-        1579 000121-33-5  1
                 HYDROXY-3-METHOXY

21  57.761  1.22 C:\Database\DEMO1.L\Adam.L
                 19.94 PHENYL ETHANE<1-NITRO-2->           1197 000000-00-0  9
                 42.03 PHENYL ETHYL OCTANOATE              1330 005457-70-5  9
                 34.46 PHENYL ETHYL HEXANOATE              1328 006290-37-5  9

22  68.463  1.71 C:\Database\DEMO1.L\Adam.L
                 11.07 METHYL BENZOATE                      983 000093-58-3  9
                 32.37 HEXENYL BENZOATE<2E->                295 076841-70-8  9
                 21.28 PHENYL ETHYL OXYACETALDEHYDE        1331 041847-88-5  9
                   (=ACETALDEHYDE,
```

Library Search Report

Data Path : D:\data\New Folder\QAL feb.2016\
Data File : ROSE ABSOLUTE XXXI-58.D
Acq On : 9 May 2022 15:32
Operator :
Sample : ROSE ABSOLUTE XXXI-58
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\NIST05.L Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: ChemStation Integrator - autoint1.e

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
1	2.210	0.54	C:\Database\NIST05.L			
			Ethyl alcohol	95	000064-17-5	74
			Ethyl alcohol	93	000064-17-5	64
			Ethyl alcohol	94	000064-17-5	43
2	18.438	51.61	C:\Database\NIST05.L			
			Phenylethyl Alcohol	9611	000060-12-8	94
			Phenylethyl Alcohol	9613	000060-12-8	91
			Phenylethyl Alcohol	9612	000060-12-8	91
3	22.305	4.27	C:\Database\NIST05.L			
			6-Octen-1-ol, 3,7-dimethyl-, (R)-	27134	001117-61-9	89
			6-Octen-1-ol, 3,7-dimethyl-	27110	000106-22-9	50
			cis-2,6-Dimethyl-2,6-octadiene	16346	002492-22-0	43
4	23.235	6.43	C:\Database\NIST05.L			
			2,6-Octadien-1-ol, 3,7-dimethyl-	25634	000624-15-7	64
			.beta.-Myrcene	15177	000123-35-3	60
			2,6-Octadien-1-ol, 3,7-dimethyl-, acetate, (Z)-	54283	000141-12-8	59
5	23.310	1.32	C:\Database\NIST05.L			
			Acetic acid, 2-phenylethyl ester	31818	000103-45-7	90
			Acetic acid, 2-phenylethyl ester	31806	000103-45-7	90
			Bromoacetic acid, 2-phenylethyl es ter	86179	003785-33-9	83
6	27.327	0.61	C:\Database\NIST05.L			
			Propanoic acid, 2-methyl-, 3,7-dim ethyl-2,6-octadienyl ester, (E)-	74356	002345-26-8	86
			2,6-Octadien-1-ol, 3,7-dimethyl-, (Z)-	25690	000106-25-2	83
			Geranyl bromide	68113	035719-26-7	80
7	29.365	0.36	C:\Database\NIST05.L			
			Imidazole-5-butyric acid, 4,.gamma .,.gamma.-trimethyl-	54063	1000129-14-8	43
			4-(2,6,6-Trimethyl-cyclohex-1-enyl)-butan-2-ol	54448	1000194-68-8	38
			Bromo-4-fluoroacetophenone	68800	000403-29-2	38
8	30.520	0.74	C:\Database\NIST05.L			
			1,6-Cyclodecadiene, 1-methyl-5-met hylene-8-(1-methylethyl)-, [s-(E,E)]-	59960	023986-74-5	97
			1H-Cyclopenta[1,3]cyclopropa[1,2]b enzene, octahydro-7-methyl-3-methy lene-4-(1-methylethyl)-, [3aS-(3a. alpha.,3b.beta.,4.beta.,7.alpha.,7 aS*)]-	60104	013744-15-5	97
			1H-Cyclopenta[1,3]cyclopropa[1,2]b enzene, octahydro-7-methyl-3-methy lene-4-(1-methylethyl)-, [3aS-(3a. alpha.,3b.beta.,4.beta.,7.alpha.,7 aS*)]-	60103	013744-15-5	93

9	40.592	5.82	C:\Database\NIST05.L			
			9-Eicosene, (E)-	112106	074685-29-3	94
			1-Nonadecene	102860	018435-45-5	94
			1-Pentadecene	64457	013360-61-7	91
10	40.993	0.86	C:\Database\NIST05.L			
			Nonadecane	104271	000629-92-5	99
			Nonadecane	104272	000629-92-5	97
			Eicosane	113490	000112-95-8	91
11	42.068	1.36	C:\Database\NIST05.L			
			n-Hexadecanoic acid	96235	000057-10-3	98
			n-Hexadecanoic acid	96233	000057-10-3	98
			n-Hexadecanoic acid	96234	000057-10-3	95
12	42.191	0.25	C:\Database\NIST05.L			
			9-Eicosene, (E)-	112106	074685-29-3	97
			3-Eicosene, (E)-	112107	074685-33-9	97
			5-Eicosene, (E)-	112105	074685-30-6	93
13	43.582	0.38	C:\Database\NIST05.L			
			3-Eicosene, (E)-	112107	074685-33-9	91
			9-Eicosene, (E)-	112106	074685-29-3	91
			Trifluoroacetic acid, n-heptadecyl ester	138900	1000216-79-2	91
14	43.785	0.39	C:\Database\NIST05.L			
			10-Heneicosene (c,t)	121168	095008-11-0	96
			Trichloroacetic acid, hexadecyl ester	166990	074339-54-1	93
			17-Pentatriacontene	183898	006971-40-0	91
15	43.897	0.37	C:\Database\NIST05.L			
			Tritetracontane	188549	007098-21-7	83
			Tetratetracontane	188837	007098-22-8	83
			Heptacosane, 1-chloro-	174384	062016-79-9	74
16	44.411	0.11	C:\Database\NIST05.L			
			8-Hexadecyne	73060	019781-86-3	58
			9-Eicosyne	110846	071899-38-2	58
			Bicyclo[2.2.2]octane, 2-methyl-	10368	000766-53-0	53
17	44.507	0.26	C:\Database\NIST05.L			
			1,4-Cyclooctadiene, (Z,Z)-	5311	016327-22-3	86
			9,12,15-Octadecatrienoic acid, (Z, Z,Z)-	110773	000463-40-1	86
			9,12,15-Octadecatrien-1-ol, (Z,Z,Z)-	101506	000506-44-5	83
18	44.769	0.29	C:\Database\NIST05.L			
			Octadecanoic acid	114820	000057-11-4	91
			Octadecanoic acid	114822	000057-11-4	78
			Octadecanoic acid	114821	000057-11-4	70
19	46.716	1.36	C:\Database\NIST05.L			
			9-Tricosene, (Z)-	138119	027519-02-4	99
			9-Tricosene, (Z)-	138118	027519-02-4	99
			Trichloroacetic acid, hexadecyl ester	166990	074339-54-1	94
20	57.247	6.18	C:\Database\NIST05.L			
			.gamma.-Tocopherol	174833	007616-22-0	96
			.gamma.-Tocopherol	174832	007616-22-0	94
			.beta.-Tocopherol	174831	000148-03-8	74
21	57.761	1.22	C:\Database\NIST05.L			

			Oxalic acid, hexadecyl 2-phenyleth yl ester	175188	1000309-67-0 87
			Oxalic acid, dodecyl 2-phenylethyl ester	158095	1000309-66-6 86
			Oxalic acid, 2-phenylethyl tridecy l ester	163893	1000309-66-7 86
22	68.463	1.71	C:\Database\NIST05.L		
			Benzene, 1,1'-(1,2-dimethyl-1,2-et hanediyl)bis-, (R*,R*)-(.+/-.)-	64500	002726-21-8 47
			Benzene, 1,1'-(1,2-dimethyl-1,2-et hanediyl)bis-, (R*,S*)-	64498	004613-11-0 40
			Benzene, 1,1'-(1,2-dimethyl-1,2-et hanediyl)bis-	64496	005789-35-5 38

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Library Search Report

Data Path : D:\data\New Folder\QAL feb.2016\
Data File : ROSE ABSOLUTE XXXI-58.D
Acq On : 9 May 2022 15:32
Operator :
Sample : ROSE ABSOLUTE XXXI-58
Misc :
ALS Vial : 1 Sample Multiplier: 1

Search Libraries: C:\Database\Flavor2.L Minimum Quality: 0

Unknown Spectrum: Apex
Integration Events: ChemStation Integrator - autoint1.e

Pk#	RT	Area%	Library/ID	Ref#	CAS#	Qual
1	2.210	0.54	C:\Database\Flavor2.L			
			Ethyl alcohol	26	000064-17-5	86
			Isopropyl formate	8	000625-55-8	3
			2-Heptanol	180	000543-49-7	2
2	18.438	51.61	C:\Database\Flavor2.L			
			Phenethyl alcohol	318	000060-12-8	95
			Phenethylamine	115	000064-04-0	2
			alpha,alpha-Dimethylphenethyl alcohol	327	000100-86-7	2
3	22.305	4.27	C:\Database\Flavor2.L			
			Geraniol	283	000106-24-1	55
			Geranyl formate	329	000105-86-2	38
			Farnesol (E,E-)	388	000106-28-5	25
4	23.235	6.43	C:\Database\Flavor2.L			
			Geraniol	283	000106-24-1	72
			Farnesol (E,E-)	388	000106-28-5	40
			Geranyl acetate	399	000105-87-3	38
5	23.310	1.32	C:\Database\Flavor2.L			
			Phenethyl acetate	293	000103-45-7	90
			xxx	291	000000-00-0	86
			Phenethyl formate	414	000104-62-1	78
6	27.327	0.61	C:\Database\Flavor2.L			
			Geranyl formate	329	000105-86-2	80
			Nerol, cis-geraniol	253	000106-25-2	74
			Geraniol	283	000106-24-1	56
7	29.365	0.36	C:\Database\Flavor2.L			
			No matches found			
8	30.520	0.74	C:\Database\Flavor2.L			
			Valencene	156	004630-07-3	12
9	33.708	13.56	C:\Database\Flavor2.L			
			Piperonal; heliotropin	67	000120-57-0	4
10	40.592	5.82	C:\Database\Flavor2.L			
			Lauryl alcohol	90	000112-53-8	38
			trans-2-Heptenal	207	018829-55-5	37
			Decyl acetate	266	000112-17-4	32
11	40.993	0.86	C:\Database\Flavor2.L			
			n-Tetracosane	412	000646-31-1	91
			n-Tetradecane	408	000629-59-4	78
			n-Eicosane	411	000112-95-8	74
12	42.068	1.36	C:\Database\Flavor2.L			
			Palmitic acid	61	000057-10-3	87
			Stearic acid	389	000057-11-4	58
			Myristic acid	60	000544-63-8	42

13	42.191	0.25	C:\Database\Flavor2.L			
			Lauryl alcohol	90	000112-53-8	47
			trans-2-Heptenal	207	018829-55-5	25
			trans-2-Hexenal	189	006728-26-3	25
14	43.582	0.38	C:\Database\Flavor2.L			
			Lauryl alcohol	90	000112-53-8	38
			trans-2-Heptenal	207	018829-55-5	32
			trans-2-Hexenal	189	006728-26-3	32
15	43.785	0.39	C:\Database\Flavor2.L			
			Lauryl alcohol	90	000112-53-8	37
			1-Decanol	86	000112-30-1	27
			Lauric aldehyde	89	000112-54-9	16
16	43.897	0.37	C:\Database\Flavor2.L			
			n-Tetracosane	412	000646-31-1	52
			n-Eicosane	411	000112-95-8	52
			n-Octadecane	410	000593-45-3	50
17	44.411	0.11	C:\Database\Flavor2.L			
			omega-6-Hexadecalactone, Ambrettol ide	71	007779-50-2	38
			cis-6-Nonen-1-ol	106	035854-86-5	22
			Citronellol	223	000106-22-9	16
18	44.507	0.26	C:\Database\Flavor2.L			
			Farnesol (E,E-)	388	000106-28-5	9
			Methyl octine carbonate	254	000111-80-8	7
			Geranyl acetate	399	000105-87-3	6
19	44.769	0.29	C:\Database\Flavor2.L			
			Stearic acid	389	000057-11-4	64
			n-Decane	405	000124-18-5	10
			Pentadecane	400	000629-62-9	10
20	46.716	1.36	C:\Database\Flavor2.L			
			Lauryl alcohol	90	000112-53-8	40
			Lauric aldehyde	89	000112-54-9	33
			1-Decanol	86	000112-30-1	32
21	57.247	6.18	C:\Database\Flavor2.L			
			Vanillin	24	000121-33-5	1
22	57.761	1.22	C:\Database\Flavor2.L			
			Phenethyloctanoate	139	005457-70-5	38
			Phenethyl isobutyrate	322	000103-48-0	9
			Phenethyl propionate	415	000122-70-3	2
23	68.463	1.71	C:\Database\Flavor2.L			
			Methyl benzoate	251	000093-58-3	2
			Benzaldehyde	359	000100-52-7	1

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