

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
 Data File : VW029523.D
 Acq On : 11 Jul 2024 12:46
 Operator : SY/MD
 Sample : P3178-03
 Misc : 4.66g/10mL/MSVOA_W/SOIL/A
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 A4BS8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
 Title : SFAM01.0

Signal : TIC: VW029523.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.363	71	78	97	rBV	71033	176345	16.25%	2.393%
2	2.899	158	166	182	rBV	51009	151260	13.94%	2.052%
3	3.283	228	229	232	rBV3	576	536	0.05%	0.007%
4	3.314	232	234	241	rVB4	580	849	0.08%	0.012%
5	3.399	241	248	256	rVB4	2290	5525	0.51%	0.075%
6	3.460	256	258	260	rBV2	526	538	0.05%	0.007%
7	3.521	266	268	270	rVB3	610	601	0.06%	0.008%
8	3.686	292	295	296	rBV3	625	590	0.05%	0.008%
9	4.021	338	350	374	rVV	124993	411281	37.91%	5.581%
10	4.393	410	411	415	rVB3	640	556	0.05%	0.008%
11	4.655	451	454	457	rBV4	393	505	0.05%	0.007%
12	4.685	457	459	462	rVB4	726	651	0.06%	0.009%
13	4.917	483	497	510	rBV2	21523	82867	7.64%	1.124%
14	5.161	535	537	539	rBV2	499	506	0.05%	0.007%
15	5.283	555	557	559	rBV2	709	660	0.06%	0.009%
16	5.350	562	568	569	rBV3	576	910	0.08%	0.012%
17	5.765	629	636	640	rBV5	451	984	0.09%	0.013%
18	5.941	656	665	677	rBV7	3920	12665	1.17%	0.172%
19	6.069	684	686	690	rVB3	543	661	0.06%	0.009%
20	6.338	727	730	732	rBV	778	587	0.05%	0.008%
21	6.508	755	758	761	rBV4	486	631	0.06%	0.009%
22	6.911	821	824	826	rBV3	631	922	0.08%	0.013%
23	7.081	844	852	859	rBV2	17325	56757	5.23%	0.770%
24	7.276	883	884	890	rVB3	638	661	0.06%	0.009%
25	7.380	899	901	904	rVB4	534	641	0.06%	0.009%
26	7.411	904	906	909	rBV3	526	870	0.08%	0.012%
27	7.648	933	945	956	rBV	188742	461241	42.51%	6.259%
28	7.776	965	966	971	rVB4	850	644	0.06%	0.009%
29	7.837	974	976	981	rVB3	409	673	0.06%	0.009%
30	8.032	1005	1008	1011	rVB2	419	622	0.06%	0.008%
31	8.276	1038	1048	1066	rBV2	373144	1084990	100.00%	14.722%
32	8.465	1075	1079	1080	rBV3	613	663	0.06%	0.009%
33	8.508	1082	1086	1089	rVB2	555	688	0.06%	0.009%
34	8.685	1112	1115	1116	rBV3	903	906	0.08%	0.012%
35	8.843	1131	1141	1154	rBV	415223	825175	76.05%	11.197%
36	9.038	1166	1173	1178	rBV2	3856	9085	0.84%	0.123%

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Integrator: RTE

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
 Title : SFAM01.0

37	9.276	1202	1212	1222	rBV	248030	512142	47.20%	6.949%
38	9.453	1237	1241	1244	rVB5	464	626	0.06%	0.008%
39	9.758	1289	1291	1297	rVB5	743	1015	0.09%	0.014%
40	9.886	1309	1312	1318	rVB6	1023	1889	0.17%	0.026%
41	10.032	1321	1336	1347	rBV	97498	187370	17.27%	2.542%
42	10.117	1347	1350	1351	rBV3	607	572	0.05%	0.008%
43	10.245	1367	1371	1375	rVB5	761	1469	0.14%	0.020%
44	10.325	1375	1384	1393	rBV	408865	724254	66.75%	9.827%
45	10.508	1410	1414	1415	rBV2	515	620	0.06%	0.008%
46	10.575	1418	1425	1434	rBV	60308	110709	10.20%	1.502%
47	10.697	1441	1445	1452	rVB	8965	16118	1.49%	0.219%
48	10.751	1452	1454	1455	rBV3	664	502	0.05%	0.007%
49	10.843	1464	1469	1475	rVB4	2814	4922	0.45%	0.067%
50	10.922	1475	1482	1501	rBV3	70506	159415	14.69%	2.163%
51	11.172	1518	1523	1530	rVV5	2732	4872	0.45%	0.066%
52	11.343	1547	1551	1554	rBV3	816	1300	0.12%	0.018%
53	11.398	1554	1560	1566	rBV4	2904	4989	0.46%	0.068%
54	11.477	1571	1573	1577	rVB4	539	732	0.07%	0.010%
55	11.520	1577	1580	1581	rBV2	517	591	0.05%	0.008%
56	11.629	1588	1598	1606	rBV	403724	666418	61.42%	9.043%
57	11.782	1618	1623	1624	rVB5	381	569	0.05%	0.008%
58	11.831	1627	1631	1632	rVV2	1521	1794	0.17%	0.024%
59	11.867	1635	1637	1642	rVB4	1611	2483	0.23%	0.034%
60	12.026	1657	1663	1666	rBV6	541	1035	0.10%	0.014%
61	12.068	1666	1670	1671	rBV3	842	1113	0.10%	0.015%
62	12.166	1683	1686	1689	rVB4	946	1081	0.10%	0.015%
63	12.245	1694	1699	1702	rVB6	846	1274	0.12%	0.017%
64	12.337	1709	1714	1716	rBV4	866	1588	0.15%	0.022%
65	12.458	1728	1734	1739	rBV6	1699	4117	0.38%	0.056%
66	12.599	1748	1757	1763	rBV2	35703	73745	6.80%	1.001%
67	12.690	1765	1772	1778	rVV	107604	181954	16.77%	2.469%
68	12.751	1778	1782	1792	rVB3	9323	22316	2.06%	0.303%
69	12.861	1797	1800	1801	rBV2	711	796	0.07%	0.011%
70	12.958	1807	1816	1821	rBV3	8881	21826	2.01%	0.296%
71	13.007	1822	1824	1827	rVV3	2838	4074	0.38%	0.055%
72	13.080	1831	1836	1839	rBV5	3330	5690	0.52%	0.077%
73	13.129	1841	1844	1845	rVV3	2637	3342	0.31%	0.045%
74	13.190	1845	1854	1861	rVB	55302	98661	9.09%	1.339%
75	13.294	1866	1871	1879	rVB	23198	38599	3.56%	0.524%
76	13.403	1879	1889	1894	rBV	81187	150818	13.90%	2.046%

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 Title : SFAM01.0

77	13.470	1897	1900	1902	rBV3	6886	9582	0.88%	0.130%
78	13.513	1902	1907	1909	rVV3	19276	37085	3.42%	0.503%
79	13.556	1909	1914	1917	rVV	191242	311544	28.71%	4.227%
80	13.598	1917	1921	1935	rVB	113705	250135	23.05%	3.394%
81	13.757	1941	1947	1955	rVB2	43630	83920	7.73%	1.139%
82	13.848	1955	1962	1967	rBV	121289	209295	19.29%	2.840%
83	13.897	1967	1970	1974	rVB	13542	16723	1.54%	0.227%
84	14.062	1992	1997	2001	rBV2	25083	38092	3.51%	0.517%
85	14.318	2034	2039	2047	rVB3	21075	33315	3.07%	0.452%
86	14.495	2063	2068	2069	rBV3	4763	6869	0.63%	0.093%
87	14.617	2085	2088	2091	rBV4	946	1497	0.14%	0.020%
88	14.684	2095	2099	2101	rVV3	4860	7272	0.67%	0.099%
89	14.714	2101	2104	2109	rVB2	7185	10420	0.96%	0.141%
90	14.830	2120	2123	2128	rBV7	782	1475	0.14%	0.020%
91	15.019	2152	2154	2161	rVB7	948	1485	0.14%	0.020%
92	15.092	2164	2166	2170	rBV4	1093	1294	0.12%	0.018%
93	15.397	2211	2216	2221	rVB3	6263	9393	0.87%	0.127%
94	15.482	2224	2230	2235	rBV	4596	10656	0.98%	0.145%
95	15.671	2258	2261	2266	rBV6	1972	3621	0.33%	0.049%
96	15.757	2272	2275	2276	rBV3	1562	1825	0.17%	0.025%
97	15.787	2277	2280	2285	rVB4	5158	7263	0.67%	0.099%
98	15.946	2302	2306	2307	rBV4	1686	2528	0.23%	0.034%
99	16.208	2345	2349	2355	rVB7	1049	2013	0.19%	0.027%
100	16.446	2386	2388	2389	rBV	911	749	0.07%	0.010%

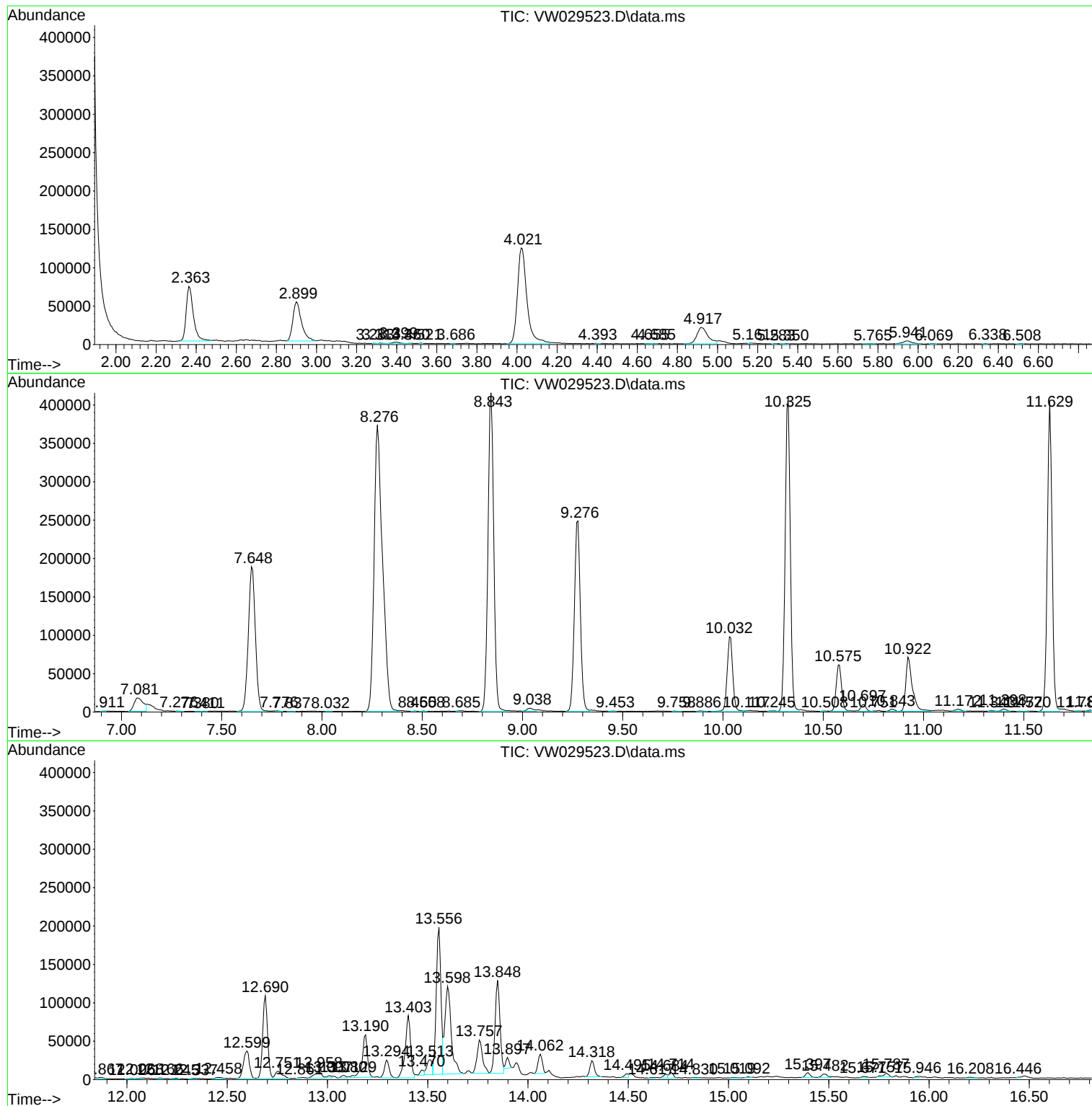
Sum of corrected areas: 7369697

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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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Instrument :
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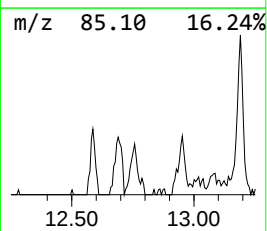
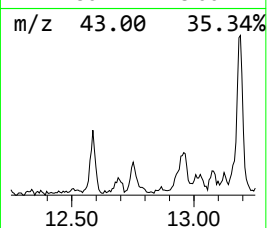
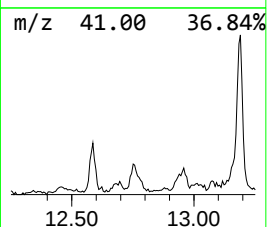
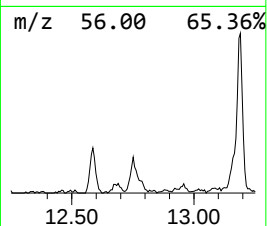
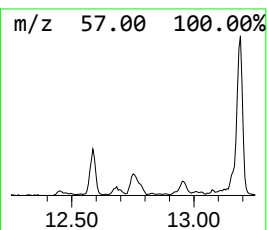
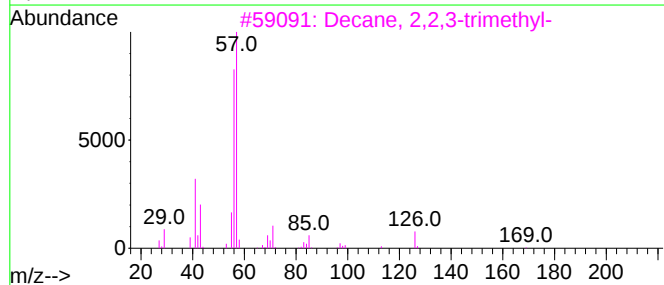
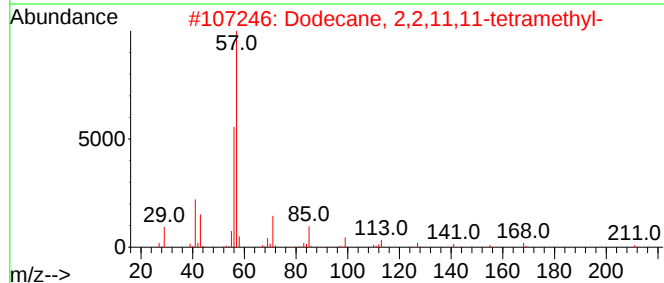
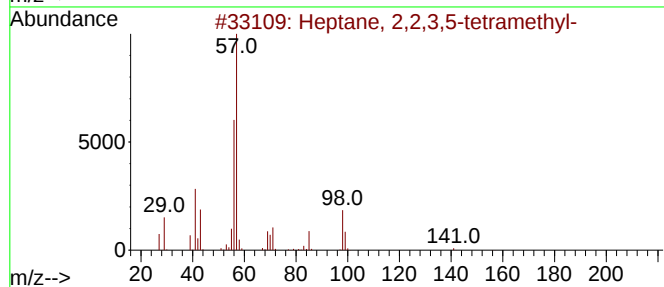
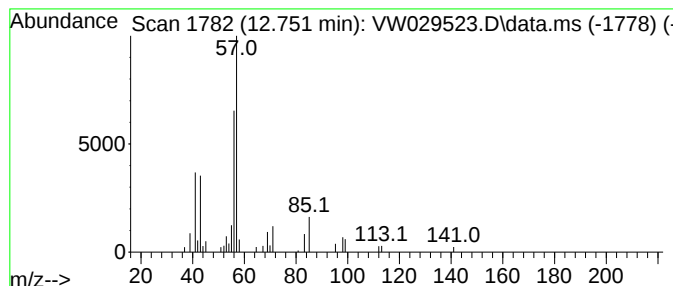
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.751	1.79 ug/L	22316	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,2,3,5-tetramethyl-	156	C11H24	061868-42-6	72
2			Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	64
3			Decane, 2,2,3-trimethyl-	184	C13H28	062338-09-4	64
4			Tetradecane, 2,2-dimethyl-	226	C16H34	059222-86-5	64
5			Octane, 2,2-dimethyl-	142	C10H22	015869-87-1	53



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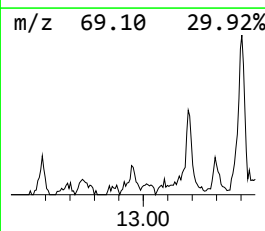
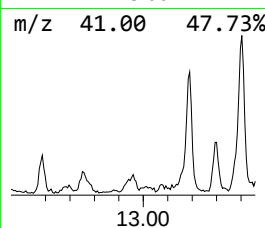
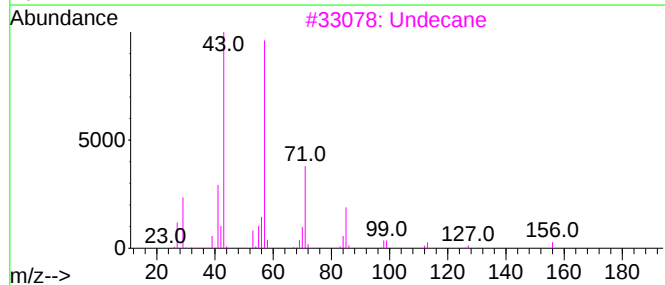
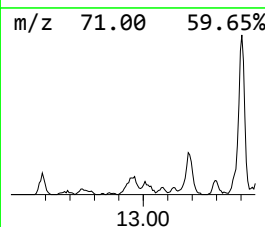
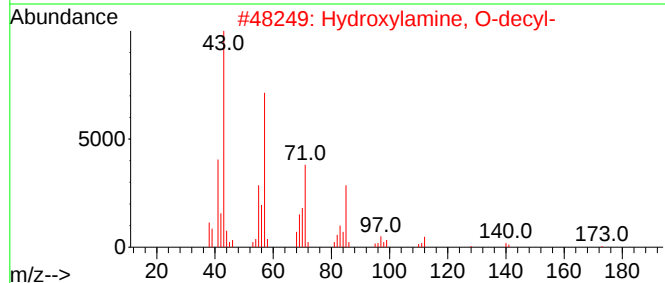
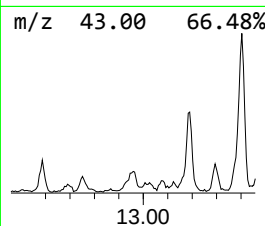
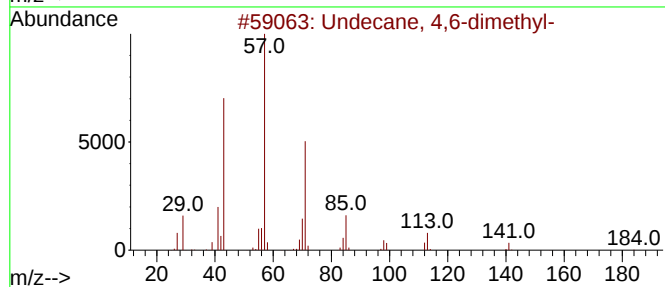
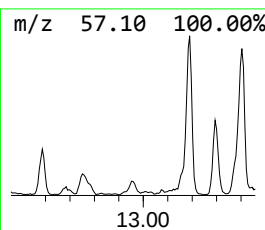
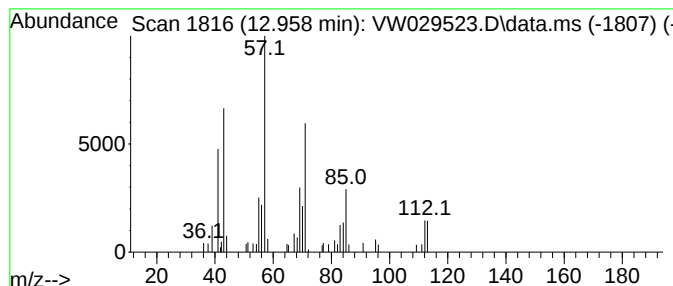
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.958	1.75 ug/L	21826	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	72
2			Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	72
3			Undecane	156	C11H24	001120-21-4	64
4			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	64
5			Octane, 2,3,3-trimethyl-	156	C11H24	062016-30-2	59



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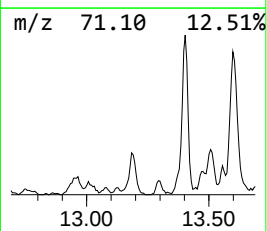
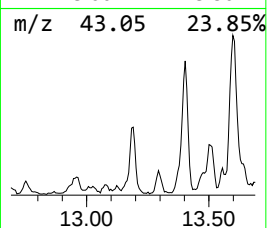
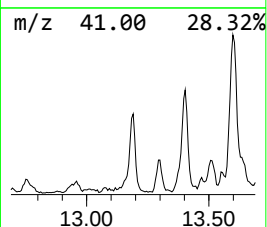
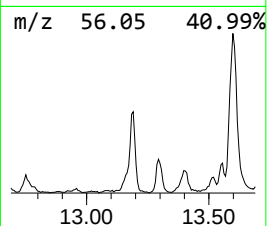
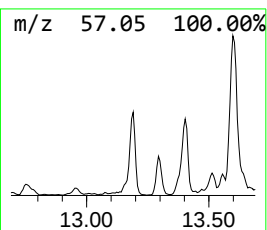
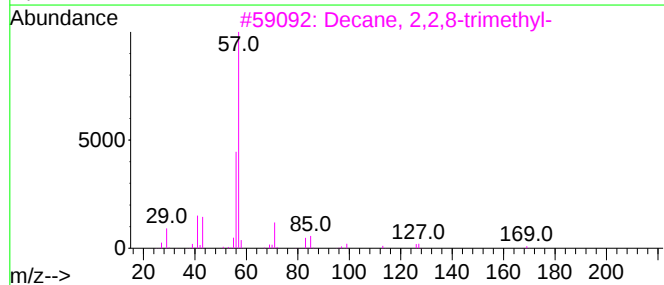
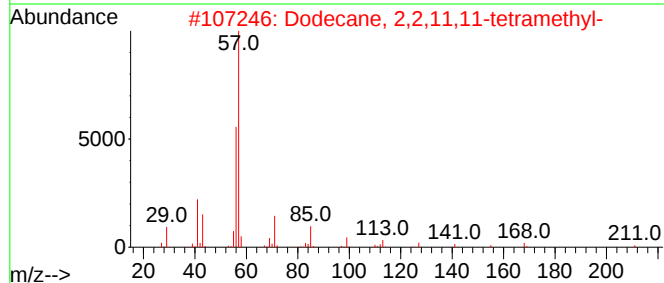
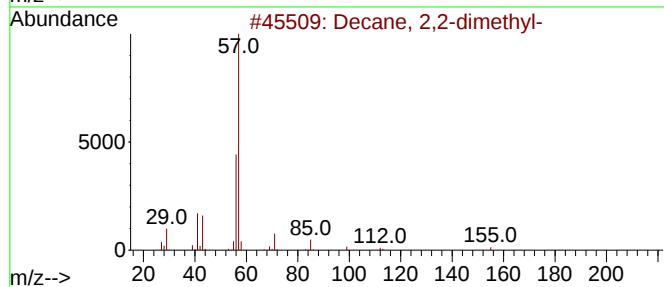
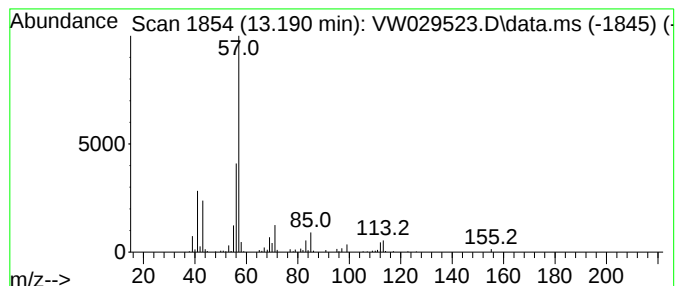
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.190	7.92 ug/L	98661	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 2,2-dimethyl-	170	C12H26	017302-37-3	78
2			Dodecane, 2,2,11,11-tetramethyl-	226	C16H34	127204-12-0	72
3			Decane, 2,2,8-trimethyl-	184	C13H28	062238-01-1	72
4			Decane, 2,2,7-trimethyl-	184	C13H28	062237-99-4	64
5			Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
Data File : VW029523.D
Acq On : 11 Jul 2024 12:46
Operator : SY/MD
Sample : P3178-03
Misc : 4.66g/10mL/MSVOA_W/SOIL/A
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A4BS8

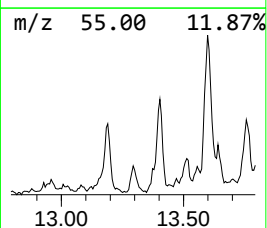
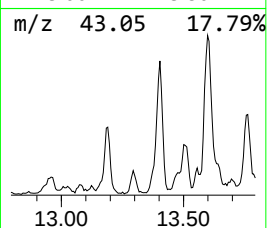
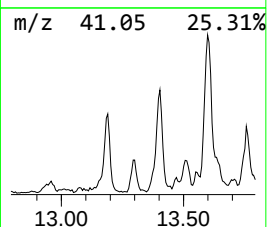
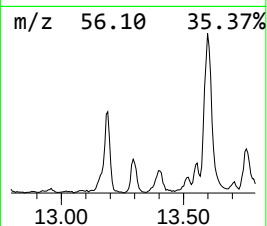
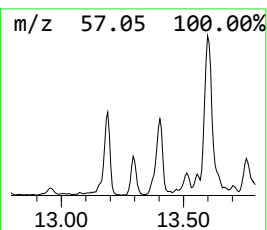
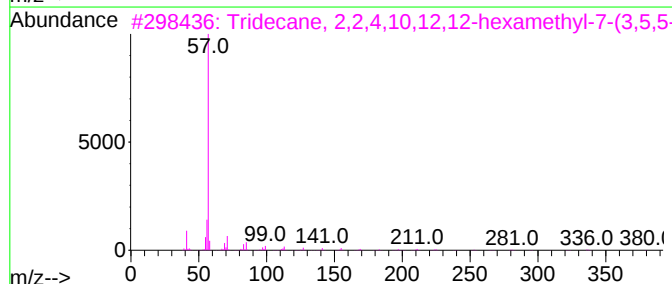
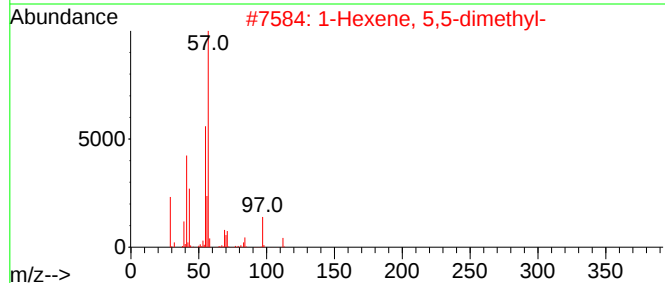
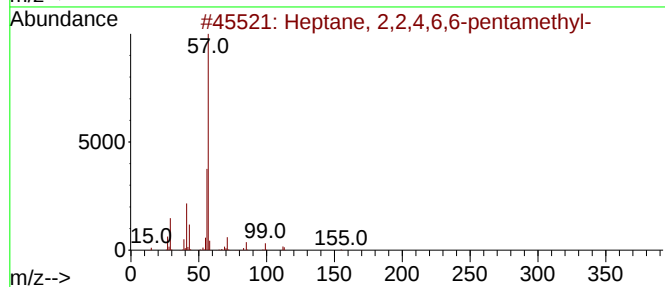
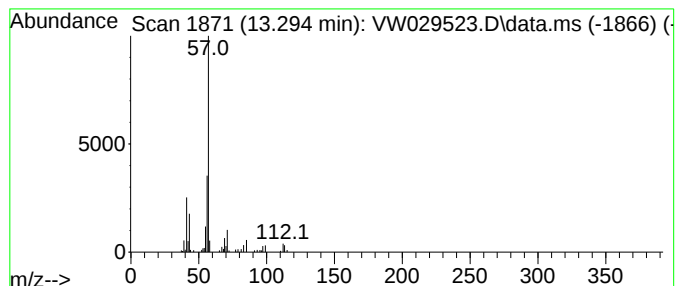
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.294	3.10 ug/L	38599	1,4-Dichlorobenzene-d4	13.556

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	78
2		1-Hexene, 5,5-dimethyl-	112	C8H16	007116-86-1	64
3		Tridecane, 2,2,4,10,12,12-hexame...	394	C28H58	003035-75-4	64
4		Heptane, 2,2,4,6,6-pentamethyl-	170	C12H26	013475-82-6	64
5		Octadecane, 2,2,4,15,17,17-hexam...	591	C42H86	055470-97-8	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
Data File : VW029523.D
Acq On : 11 Jul 2024 12:46
Operator : SY/MD
Sample : P3178-03
Misc : 4.66g/10mL/MSVOA_W/SOIL/A
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A4BS8

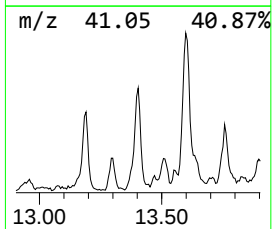
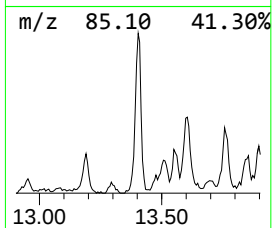
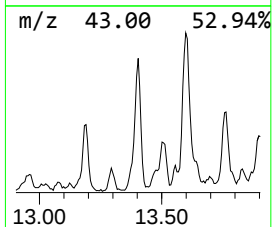
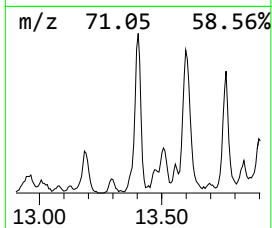
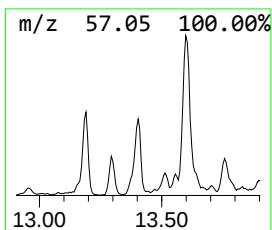
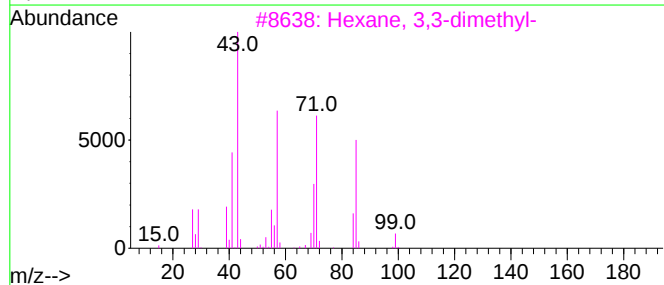
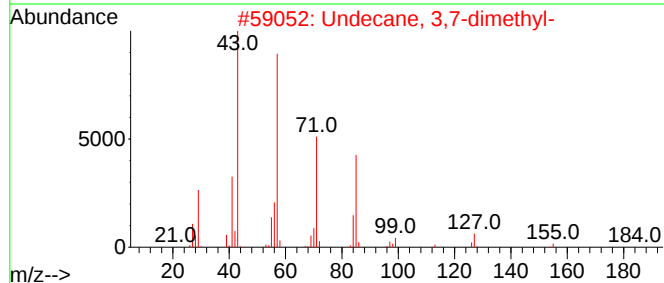
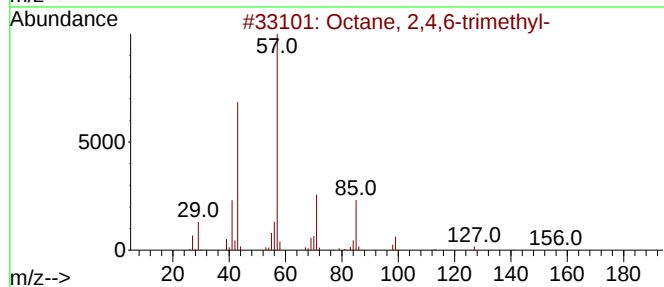
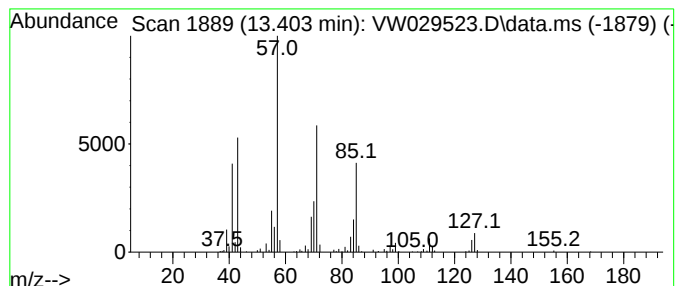
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.403	12.10 ug/L	150818	1,4-Dichlorobenzene-d4	13.556

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octane, 2,4,6-trimethyl-	156	C11H24	062016-37-9	72
2		Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	64
3		Hexane, 3,3-dimethyl-	114	C8H18	000563-16-6	50
4		Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	47
5		2-Nonanol, trimethylacetate	228	C14H28O2	1000463-21-6	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
Data File : VW029523.D
Acq On : 11 Jul 2024 12:46
Operator : SY/MD
Sample : P3178-03
Misc : 4.66g/10mL/MSVOA_W/SOIL/A
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A4BS8

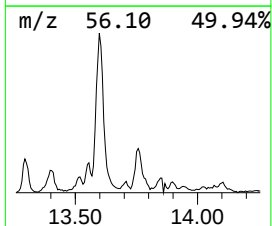
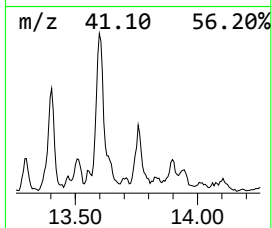
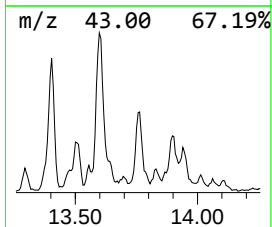
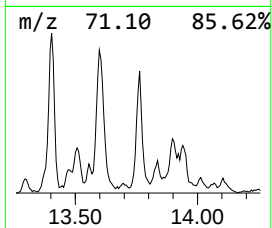
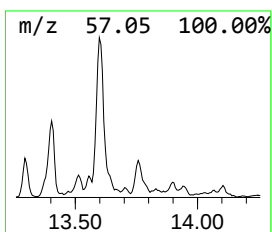
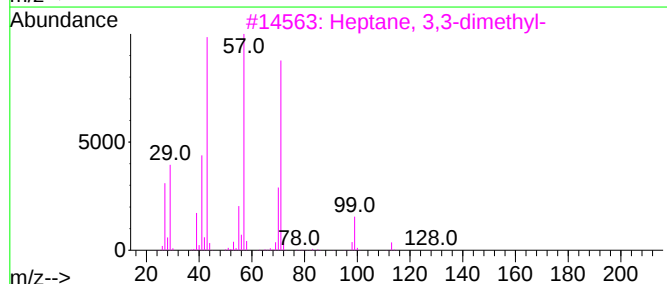
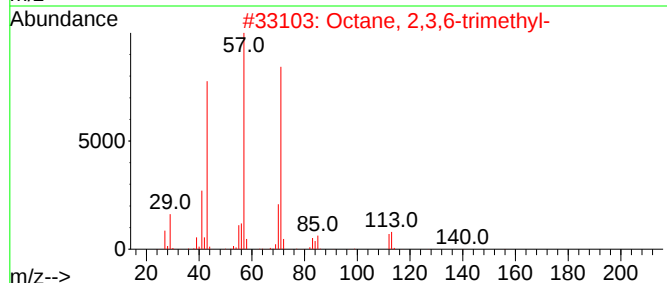
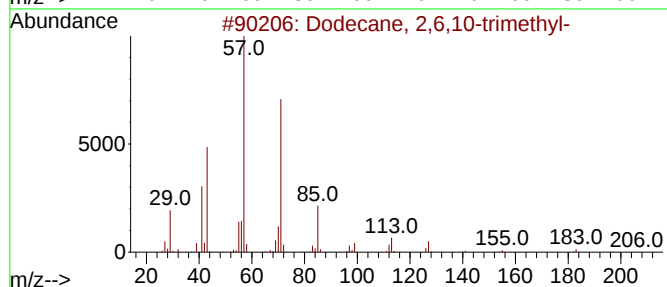
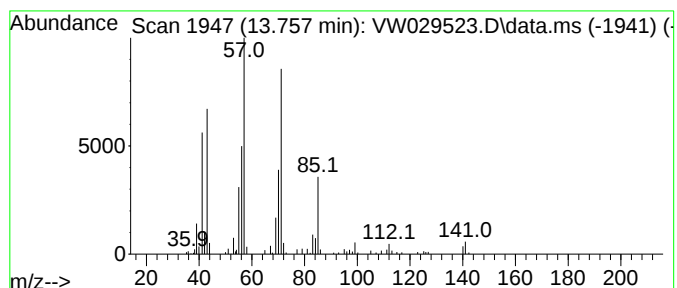
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	6.73 ug/L	83920	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	50
2			Octane, 2,3,6-trimethyl-	156	C11H24	062016-33-5	50
3			Heptane, 3,3-dimethyl-	128	C9H20	004032-86-4	50
4			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	47
5			Nonane, 2,3-dimethyl-	156	C11H24	002884-06-2	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
Data File : VW029523.D
Acq On : 11 Jul 2024 12:46
Operator : SY/MD
Sample : P3178-03
Misc : 4.66g/10mL/MSVOA_W/SOIL/A
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A4BS8

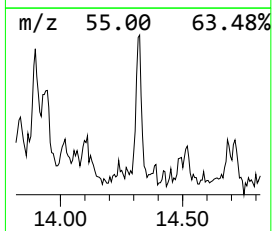
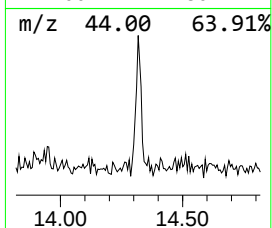
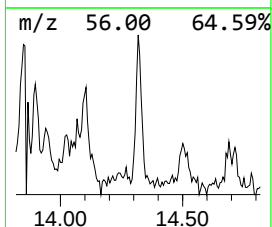
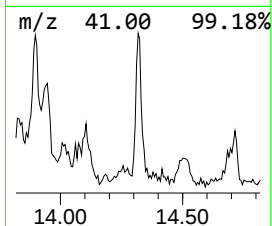
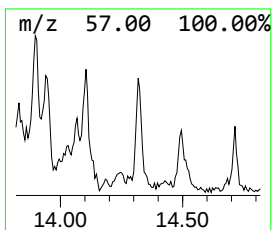
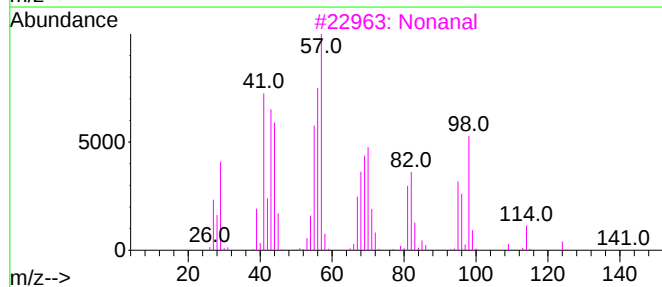
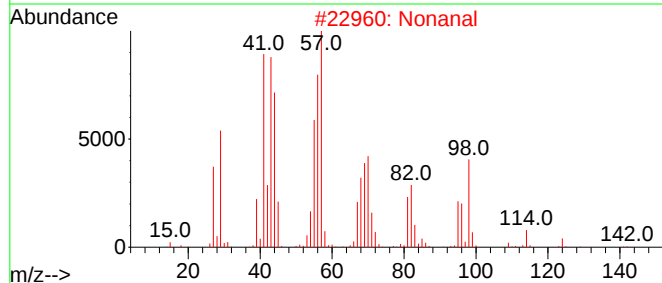
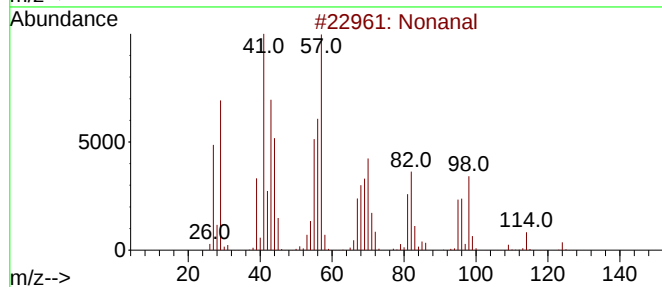
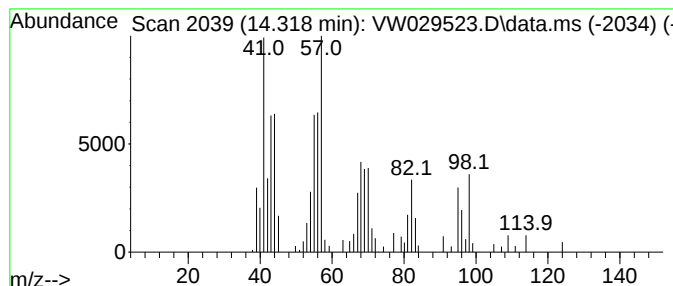
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Nonanal Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.318	2.67 ug/L	33315	1,4-Dichlorobenzene-d4	13.556

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonanal			142	C9H18O	000124-19-6	86
2	Nonanal			142	C9H18O	000124-19-6	80
3	Nonanal			142	C9H18O	000124-19-6	72
4	2-Propenoic acid, undec-10-enyl ...			224	C14H24O2	020763-72-8	35
5	Cycloheptane			98	C7H14	000291-64-5	30



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW071124\
Data File : VW029523.D
Acq On : 11 Jul 2024 12:46
Operator : SY/MD
Sample : P3178-03
Misc : 4.66g/10mL/MSVOA_W/SOIL/A
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
A4BS8

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM070824SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

						--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc	
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(DEL) Alkane: S...	12.958	1.8	ug/L	21826	3	13.556	311544	25.0	
(DEL) Alkane: S...	13.190	7.9	ug/L	98661	3	13.556	311544	25.0	
(DEL) Alkane: S...	13.294	3.1	ug/L	38599	3	13.556	311544	25.0	
(DEL) Alkane: S...	13.403	12.1	ug/L	150818	3	13.556	311544	25.0	
(DEL) Alkane: S...	13.757	6.7	ug/L	83920	3	13.556	311544	25.0	
Nonanal	14.318	2.7	ug/L	33315	3	13.556	311544	25.0	