

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122324\
 Data File : VW031513.D
 Acq On : 23 Dec 2024 17:34
 Operator : SY/MD
 Sample : P5322-19
 Misc : 6.17g/10mL/MSVOA_W/SOIL/A
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_W
 ClientSampleId :
 YE642

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 0 % of largest Peak

Max Peaks: 100

Peak Location: TOP

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

Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM121324SMA.M
 Title : SFAM01.0

Signal : TIC: VW031513.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.162	41	45	47	rBV3	4323	6278	0.75%	0.112%
2	2.357	72	77	86	rVB	41656	78032	9.31%	1.394%
3	2.503	98	101	109	rVB	5392	10741	1.28%	0.192%
4	2.893	159	165	175	rVB2	36117	74342	8.87%	1.328%
5	3.253	222	224	226	rBV3	991	616	0.07%	0.011%
6	3.917	331	333	336	rBV3	668	916	0.11%	0.016%
7	3.948	336	338	340	rVV3	711	566	0.07%	0.010%
8	4.027	340	351	361	rVV	70048	191472	22.84%	3.419%
9	4.131	364	368	378	rVB3	3193	8795	1.05%	0.157%
10	4.301	394	396	399	rVB3	957	853	0.10%	0.015%
11	4.332	399	401	403	rBV3	774	850	0.10%	0.015%
12	4.497	426	428	432	rVB4	982	915	0.11%	0.016%
13	4.582	439	442	444	rBV3	793	877	0.10%	0.016%
14	4.747	468	469	473	rVB4	757	609	0.07%	0.011%
15	4.801	476	478	481	rVB3	797	690	0.08%	0.012%
16	4.911	493	496	497	rBV3	818	834	0.10%	0.015%
17	5.088	524	525	526	rBV	810	589	0.07%	0.011%
18	5.106	526	528	532	rVV4	1099	1263	0.15%	0.023%
19	5.167	536	538	540	rBV2	525	591	0.07%	0.011%
20	5.210	543	545	547	rVV3	809	909	0.11%	0.016%
21	5.320	561	563	566	rVB4	780	653	0.08%	0.012%
22	5.502	590	593	595	rVB3	742	724	0.09%	0.013%
23	5.539	597	599	601	rBV3	728	545	0.07%	0.010%
24	5.667	615	620	621	rBV2	598	597	0.07%	0.011%
25	5.685	621	623	625	rBV	794	590	0.07%	0.011%
26	5.716	625	628	630	rBV4	728	729	0.09%	0.013%
27	5.783	635	639	641	rBV5	661	810	0.10%	0.014%
28	5.917	657	661	662	rBV4	1242	1580	0.19%	0.028%
29	5.935	662	664	666	rVV2	618	729	0.09%	0.013%
30	5.984	671	672	675	rVV3	1331	871	0.10%	0.016%
31	6.051	681	683	686	rVB4	874	720	0.09%	0.013%
32	6.088	686	689	690	rBV3	583	590	0.07%	0.011%
33	6.137	694	697	702	rBV5	767	1264	0.15%	0.023%
34	6.185	702	705	707	rBV4	430	551	0.07%	0.010%
35	6.252	715	716	719	rVB3	1000	682	0.08%	0.012%
36	6.362	730	734	736	rBV3	518	748	0.09%	0.013%

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Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM121324SMA.M
 Title : SFAM01.0

37	6.441	743	747	748	rVB3	687	825	0.10%	0.015%
38	6.460	748	750	752	rBV3	747	767	0.09%	0.014%
39	6.563	763	767	770	rBV4	965	1333	0.16%	0.024%
40	6.667	780	784	786	rBV4	408	665	0.08%	0.012%
41	6.734	793	795	797	rBV3	552	647	0.08%	0.012%
42	6.801	801	806	809	rBV4	1028	1986	0.24%	0.035%
43	6.832	809	811	815	rVB3	535	576	0.07%	0.010%
44	6.868	815	817	818	rBV2	899	560	0.07%	0.010%
45	6.911	818	824	828	rBV7	1952	4949	0.59%	0.088%
46	6.990	834	837	840	rVB3	611	629	0.08%	0.011%
47	7.088	845	853	856	rBV	16526	39958	4.77%	0.714%
48	7.386	900	902	903	rBV	1234	709	0.08%	0.013%
49	7.405	903	905	906	rVV2	704	647	0.08%	0.012%
50	7.514	919	923	924	rBV4	723	813	0.10%	0.015%
51	7.563	929	931	933	rBV3	769	765	0.09%	0.014%
52	7.648	935	945	956	rBV	96798	243913	29.09%	4.356%
53	7.996	1000	1002	1004	rBV2	1014	719	0.09%	0.013%
54	8.075	1010	1015	1018	rVB7	655	1012	0.12%	0.018%
55	8.106	1018	1020	1021	rBV2	625	541	0.06%	0.010%
56	8.142	1024	1026	1028	rBV3	815	557	0.07%	0.010%
57	8.276	1036	1048	1070	rVV2	196512	623348	74.35%	11.132%
58	8.423	1070	1072	1074	rVB2	991	741	0.09%	0.013%
59	8.575	1095	1097	1100	rVB4	585	565	0.07%	0.010%
60	8.642	1104	1108	1109	rVV3	552	654	0.08%	0.012%
61	8.709	1116	1119	1122	rBV5	1128	1170	0.14%	0.021%
62	8.843	1132	1141	1153	rBV	205260	419017	49.98%	7.483%
63	9.051	1173	1175	1176	rBV	741	648	0.08%	0.012%
64	9.087	1178	1181	1184	rVV5	882	1283	0.15%	0.023%
65	9.154	1190	1192	1193	rVV2	932	632	0.08%	0.011%
66	9.173	1193	1195	1197	rVB3	1340	972	0.12%	0.017%
67	9.270	1203	1211	1221	rBV	136391	286118	34.13%	5.110%
68	9.416	1229	1235	1238	rBV5	2846	5972	0.71%	0.107%
69	9.587	1260	1263	1267	rVB4	631	798	0.10%	0.014%
70	9.624	1267	1269	1272	rBV4	491	744	0.09%	0.013%
71	9.666	1274	1276	1277	rVV2	746	562	0.07%	0.010%
72	9.685	1277	1279	1280	rVV2	724	535	0.06%	0.010%
73	9.721	1283	1285	1287	rBV3	905	836	0.10%	0.015%
74	9.752	1288	1290	1296	rVB6	1434	1911	0.23%	0.034%
75	9.886	1308	1312	1318	rVB6	3011	6035	0.72%	0.108%
76	10.032	1330	1336	1344	rBV	73089	132026	15.75%	2.358%

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 Title : SFAM01.0

77	10.209	1363	1365	1370	rBV5	1726	3200	0.38%	0.057%
78	10.325	1376	1384	1391	rBV	269902	483096	57.62%	8.627%
79	10.428	1399	1401	1402	rVB2	949	683	0.08%	0.012%
80	10.508	1410	1414	1418	rBV5	1729	2030	0.24%	0.036%
81	10.581	1419	1426	1438	rVV	42179	79501	9.48%	1.420%
82	10.703	1440	1446	1451	rVB2	18410	34294	4.09%	0.612%
83	10.922	1476	1482	1496	rBV	57925	120740	14.40%	2.156%
84	11.434	1561	1566	1572	rVB3	16004	28579	3.41%	0.510%
85	11.629	1590	1598	1610	rBV	297467	527557	62.93%	9.421%
86	11.727	1610	1614	1622	rVB2	10065	21115	2.52%	0.377%
87	11.837	1628	1632	1636	rBV7	4110	6837	0.82%	0.122%
88	12.599	1752	1757	1763	rVB2	38027	64179	7.66%	1.146%
89	12.690	1766	1772	1778	rVV	109169	181495	21.65%	3.241%
90	12.776	1783	1786	1792	rVB4	12636	21948	2.62%	0.392%
91	12.928	1807	1811	1820	rVB7	25953	53302	6.36%	0.952%
92	13.160	1845	1849	1858	rBV8	14554	34291	4.09%	0.612%
93	13.495	1900	1904	1908	rVB	43370	66162	7.89%	1.182%
94	13.556	1908	1914	1931	rVB2	355244	838351	100.00%	14.971%
95	13.849	1956	1962	1971	rVB2	266932	523394	62.43%	9.347%
96	14.202	2017	2020	2025	rVB3	29183	43098	5.14%	0.770%
97	14.373	2046	2048	2051	rBV4	10607	13907	1.66%	0.248%
98	14.416	2052	2055	2064	rVB3	30162	60852	7.26%	1.087%
99	15.123	2167	2171	2176	rBV	77873	121338	14.47%	2.167%
100	15.354	2205	2209	2216	rVB	51560	89503	10.68%	1.598%

Sum of corrected areas: 5599711

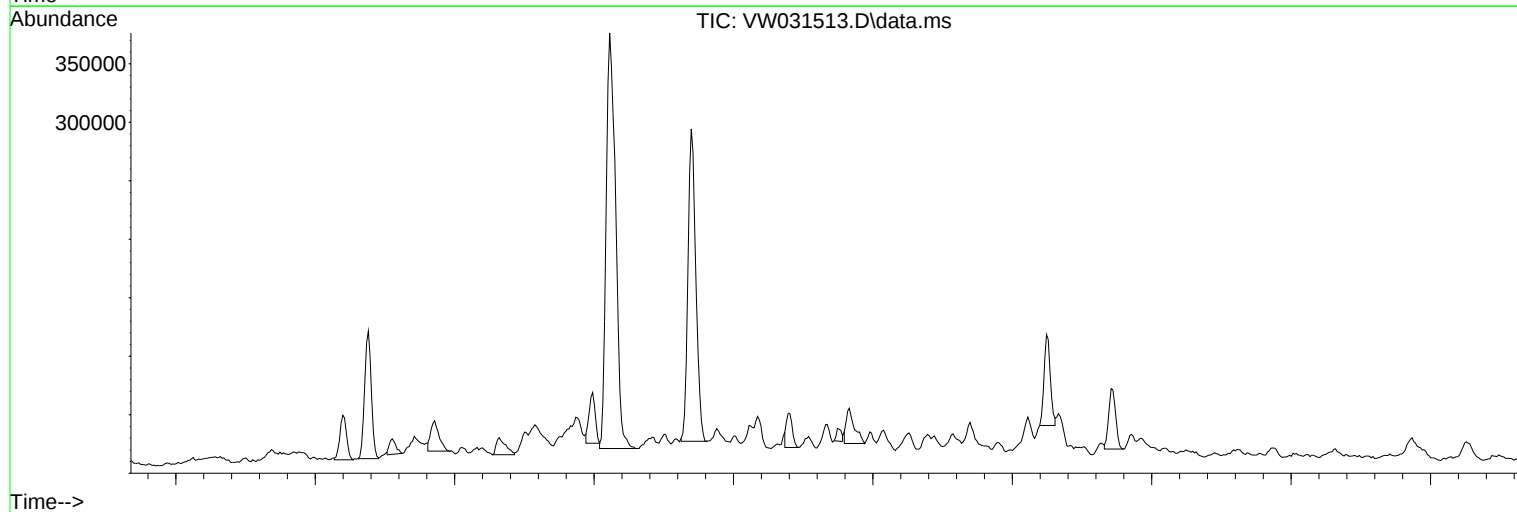
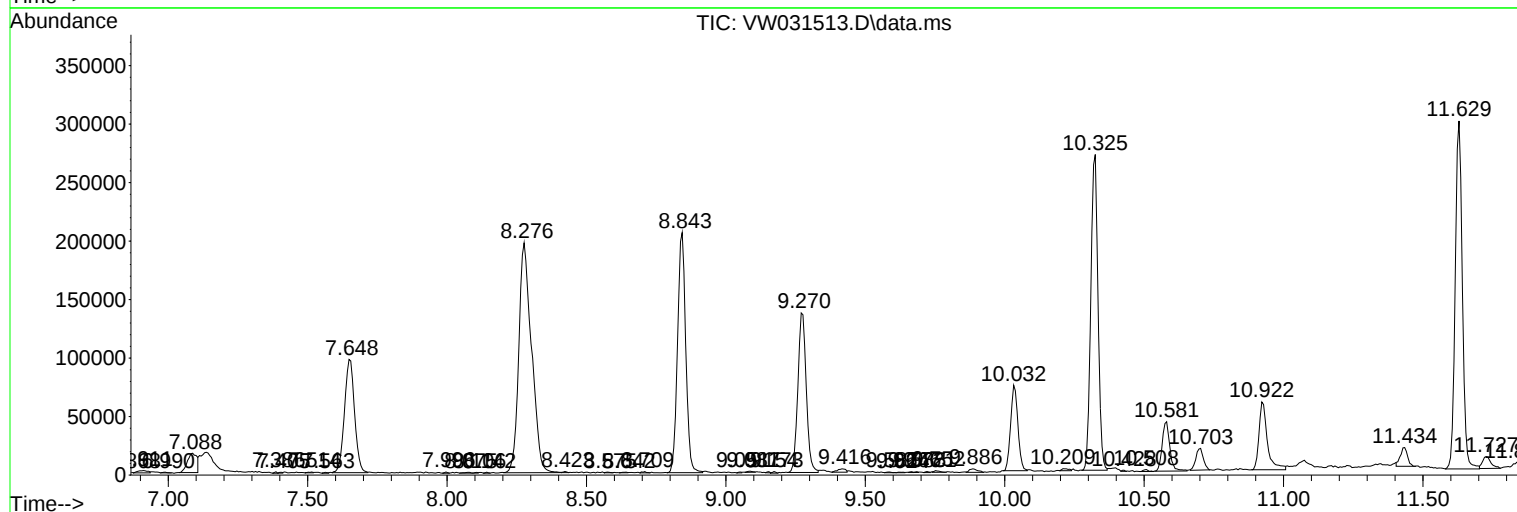
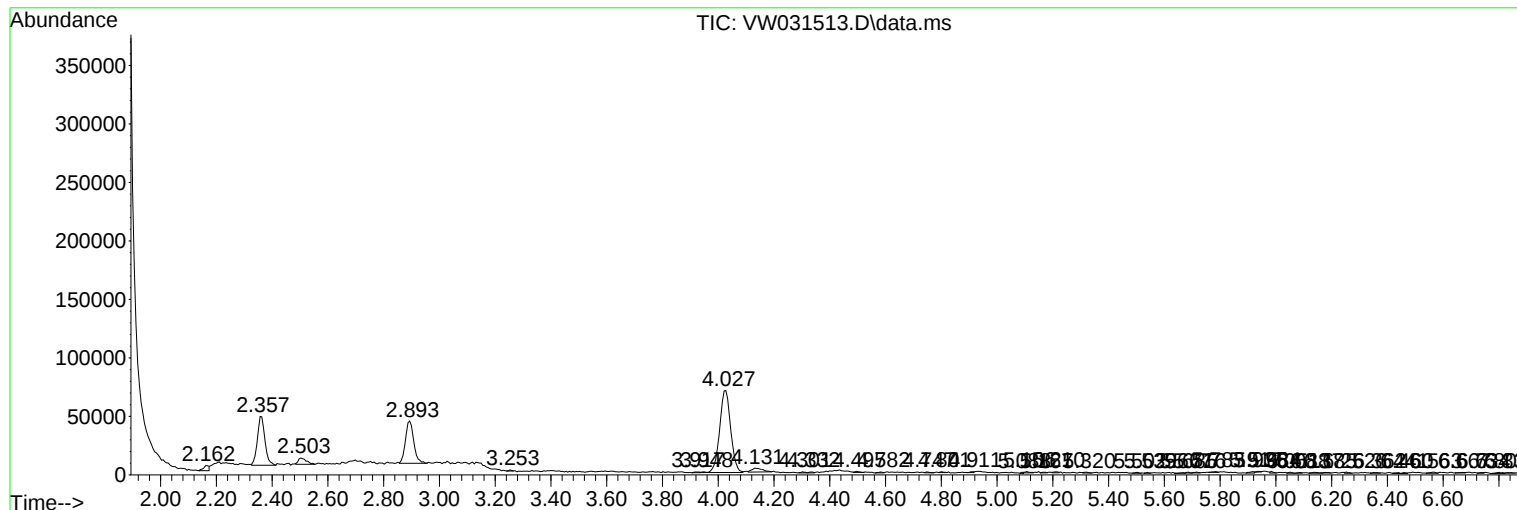
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM121324SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P



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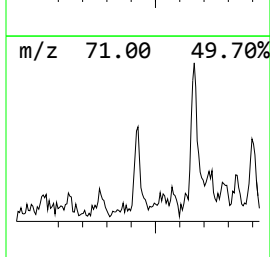
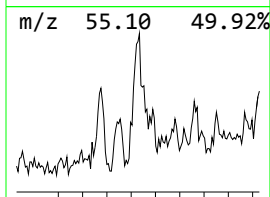
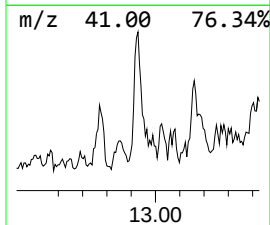
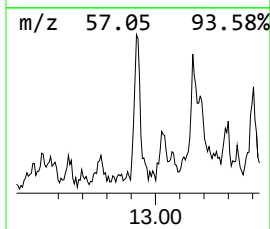
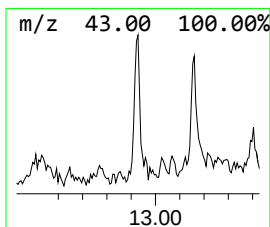
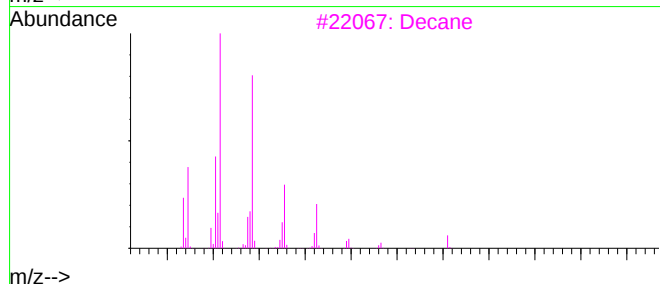
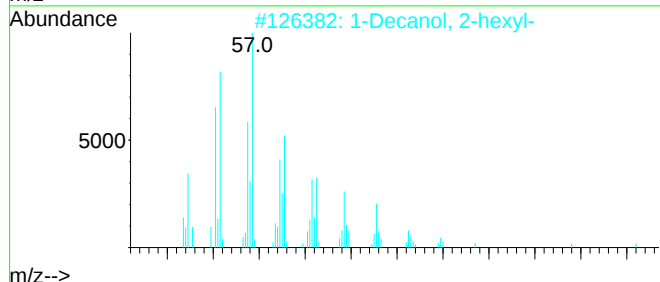
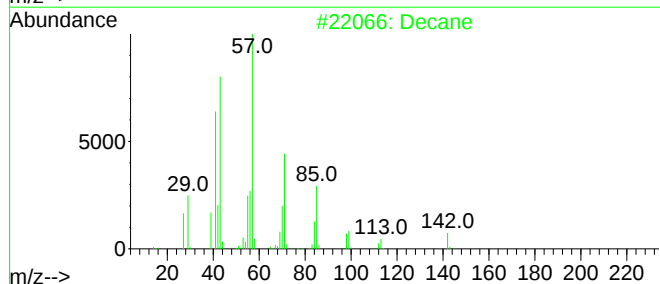
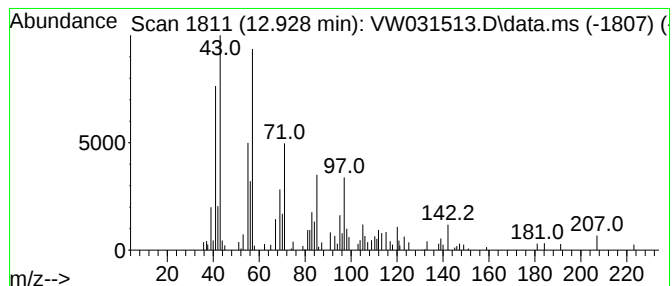
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM121324SMA.M
Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.928	1.59 ug/L	53302	1,4-Dichlorobenzene-d4	13.556	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	86
2	1-Decanol, 2-hexyl-	242	C16H34O	002425-77-6	72
3	Decane	142	C10H22	000124-18-5	62
4	Decane	142	C10H22	000124-18-5	62
5	Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	59



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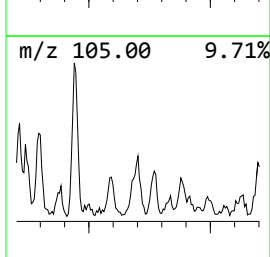
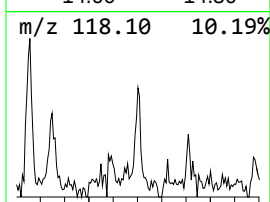
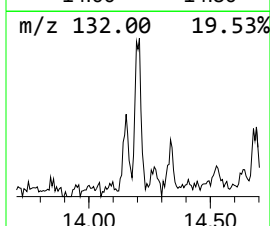
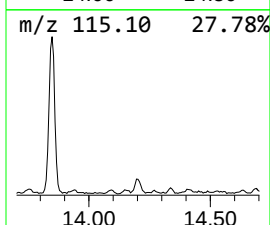
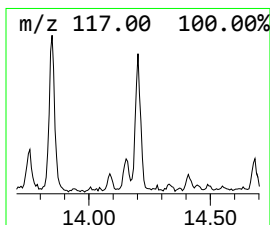
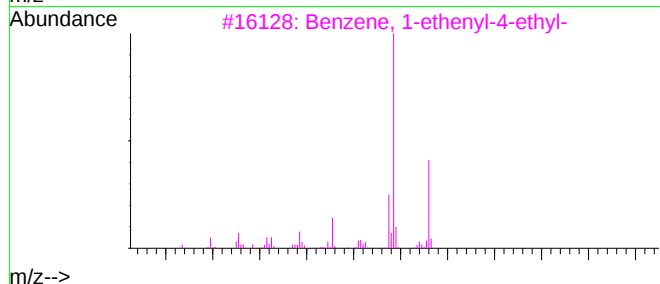
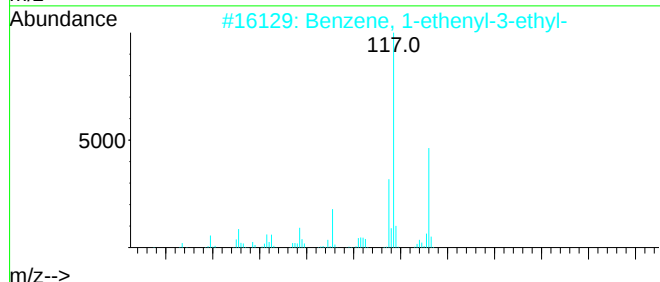
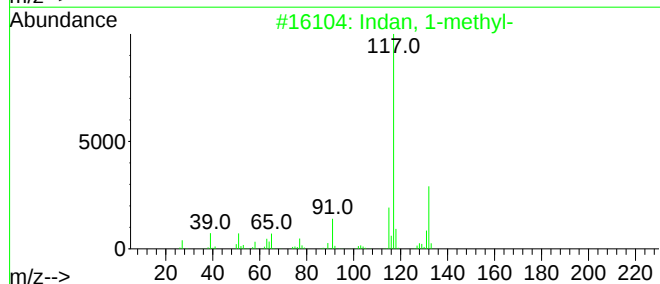
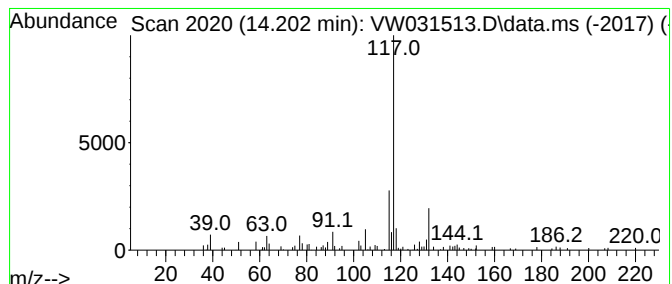
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 6 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.202	1.29 ug/L	43098	1,4-Dichlorobenzene-d4		13.556	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-		132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-		132	C10H12	007525-62-4	87
3	Benzene, 1-ethenyl-4-ethyl-		132	C10H12	003454-07-7	83
4	Benzene, 2-butenyl-		132	C10H12	001560-06-1	78
5	1-Phenyl-1-butene		132	C10H12	000824-90-8	78



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Quant Title : SFAM01.0

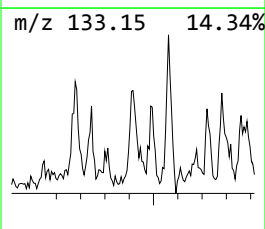
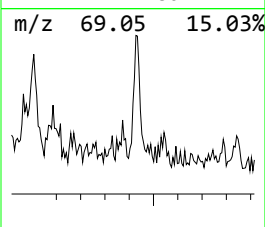
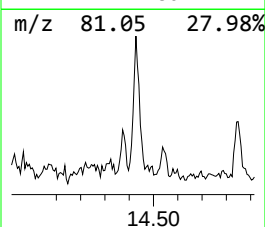
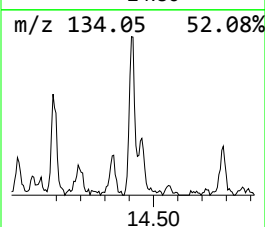
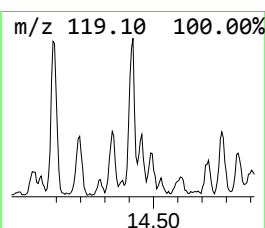
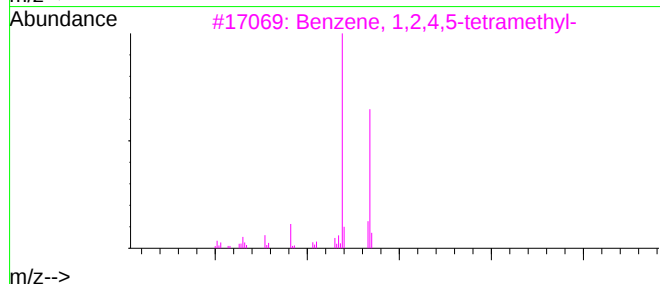
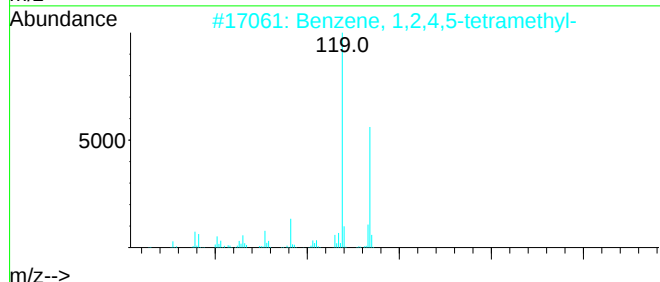
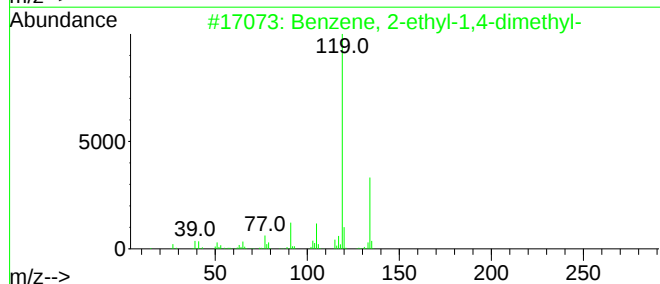
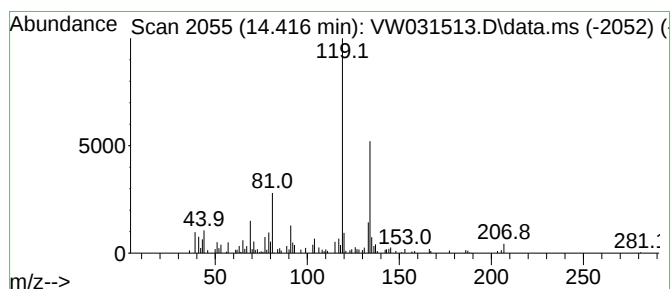
TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 7 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.416	1.81 ug/L	60852	1,4-Dichlorobenzene-d4	13.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	90
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	90
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	90
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	90
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90



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Acq On : 23 Dec 2024 17:34
Operator : SY/MD
Sample : P5322-19
Misc : 6.17g/10mL/MSVOA_W/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
YE642

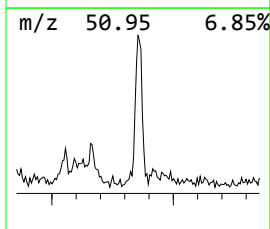
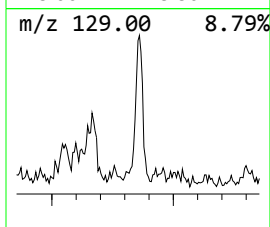
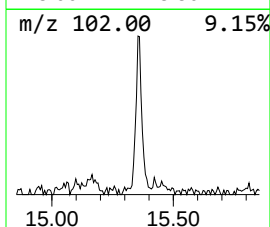
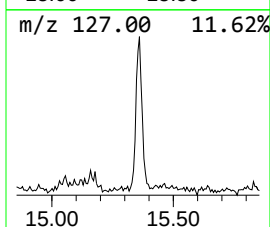
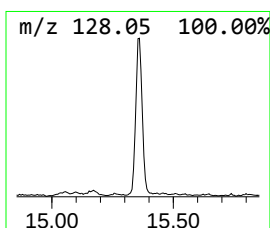
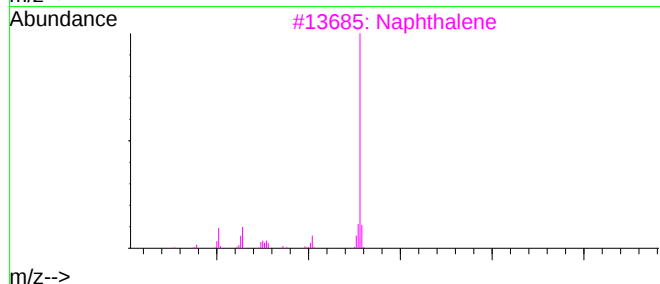
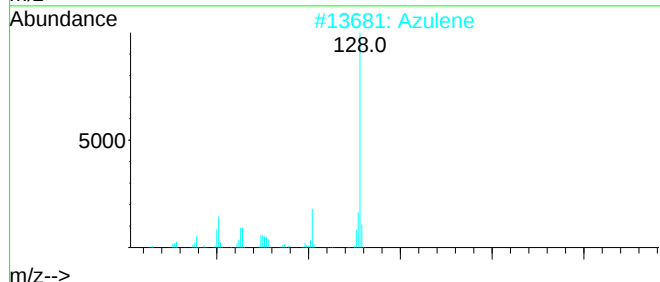
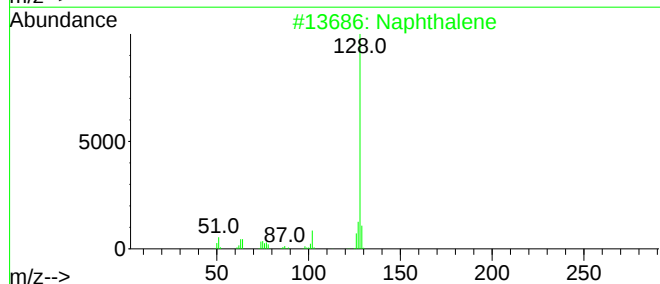
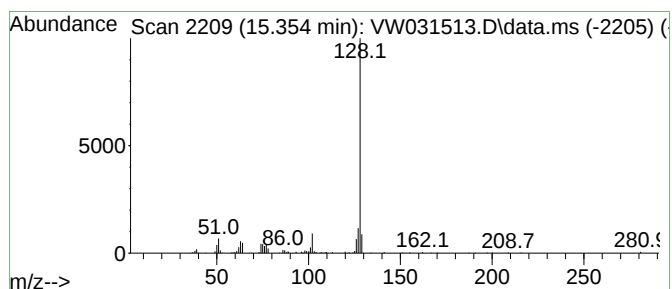
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Quant Title : SFAM01.0

TIC Library : C:\Database\NIST20.L

TIC Integration Parameters: LSCINT.P

Peak Number 8 Azulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.354	2.67 ug/L	89503	1,4-Dichlorobenzene-d4		13.556	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Naphthalene		128	C10H8	000091-20-3	95
2	Azulene		128	C10H8	000275-51-4	94
3	Naphthalene		128	C10H8	000091-20-3	90
4	Azulene		128	C10H8	000275-51-4	90
5	Azulene		128	C10H8	000275-51-4	90



Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW122324\
Data File : VW031513.D
Acq On : 23 Dec 2024 17:34
Operator : SY/MD
Sample : P5322-19
Misc : 6.17g/10mL/MSVOA_W/SOIL/A
ALS Vial : 18 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
YE642

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM121324SMA.M
Quant Title : SFAM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	12.928	1.6	ug/L	53302	3	13.556	838351	25.0
Indan, 1-methyl-	14.202	1.3	ug/L	43098	3	13.556	838351	25.0
Benzene, 2-ethy...	14.416	1.8	ug/L	60852	3	13.556	838351	25.0
Azulene	15.354	2.7	ug/L	89503	3	13.556	838351	25.0