

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
 Data File : VD076693.D
 Acq On : 03 Jul 2023 15:35
 Operator : KP/SY
 Sample : 03472-13RE
 Misc : 5.90G/10.0ml/MSVOA_D/SOIL
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 A46S5RE

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_D\Method\SFAMDLM061523SMA.M
 Title : SFAM01.0

Signal : TIC: VD076693.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.734	13	25	36	rBV	510301	964646	54.44%	9.206%
2	1.887	48	51	55	rVB2	6819	7183	0.41%	0.069%
3	2.058	76	80	81	rBV4	1846	2216	0.13%	0.021%
4	2.099	82	87	90	rBV2	2115	3534	0.20%	0.034%
5	2.140	92	94	96	rBV2	2184	1713	0.10%	0.016%
6	2.275	111	117	124	rVB2	74510	122135	6.89%	1.166%
7	2.434	141	144	147	rBV3	2922	3243	0.18%	0.031%
8	2.528	159	160	161	rBV	3409	2068	0.12%	0.020%
9	2.793	200	205	212	rBV	32713	64656	3.65%	0.617%
10	3.046	246	248	249	rVB	4032	2212	0.12%	0.021%
11	3.064	249	251	253	rBV3	2218	2143	0.12%	0.020%
12	3.693	348	358	366	rBV4	15951	42788	2.41%	0.408%
13	3.958	386	403	415	rBV5	115943	462866	26.12%	4.417%
14	4.258	446	454	460	rBV4	5325	15060	0.85%	0.144%
15	4.687	523	527	532	rBV3	1252	2112	0.12%	0.020%
16	4.752	534	538	539	rBV2	1815	2083	0.12%	0.020%
17	4.799	539	546	553	rVV4	11550	34344	1.94%	0.328%
18	4.899	560	563	566	rVB2	1384	1646	0.09%	0.016%
19	5.134	600	603	606	rVB2	1264	1632	0.09%	0.016%
20	5.311	624	633	643	rBV4	7762	24063	1.36%	0.230%
21	5.828	713	721	729	rBV6	3495	10960	0.62%	0.105%
22	6.116	758	770	780	rVV2	35220	113017	6.38%	1.079%
23	6.593	850	851	854	rBV2	1398	1550	0.09%	0.015%
24	6.622	854	856	861	rVB2	1383	2160	0.12%	0.021%
25	6.887	900	901	905	rBV2	1260	1584	0.09%	0.015%
26	6.987	908	918	922	rBV	25819	66061	3.73%	0.630%
27	7.075	923	933	950	rVB	654943	1771799	100.00%	16.908%
28	7.381	982	985	988	rBV2	1264	1837	0.10%	0.018%
29	7.569	1007	1017	1029	rVV2	79777	214879	12.13%	2.051%
30	7.793	1044	1055	1067	rBV	227956	605471	34.17%	5.778%
31	7.869	1067	1068	1072	rVV4	2510	2680	0.15%	0.026%
32	8.093	1100	1106	1107	rBV3	1503	2770	0.16%	0.026%
33	8.204	1116	1125	1140	rBV3	123612	411526	23.23%	3.927%
34	8.328	1140	1146	1152	rVV2	8649	19305	1.09%	0.184%
35	8.410	1152	1160	1171	rVB	16179	41620	2.35%	0.397%
36	8.775	1209	1222	1233	rBV	244406	492395	27.79%	4.699%

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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

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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_D\Method\SFAMDLM061523SMA.M
 Title : SFAM01.0

37	9.028	1257	1265	1273	rBV2	51626	112432	6.35%	1.073%
38	9.210	1287	1296	1303	rBV2	113692	235325	13.28%	2.246%
39	9.275	1305	1307	1312	rVV3	3750	4538	0.26%	0.043%
40	9.328	1315	1316	1319	rVB3	1892	1856	0.10%	0.018%
41	9.528	1344	1350	1351	rBV3	3237	5494	0.31%	0.052%
42	9.546	1351	1353	1355	rVV3	2235	2194	0.12%	0.021%
43	9.599	1359	1362	1365	rVB3	2207	2350	0.13%	0.022%
44	9.693	1369	1378	1387	rBV4	4202	9702	0.55%	0.093%
45	9.834	1399	1402	1403	rBV2	1878	1784	0.10%	0.017%
46	9.975	1419	1426	1433	rBV	62315	113501	6.41%	1.083%
47	10.028	1433	1435	1440	rVV5	3352	4756	0.27%	0.045%
48	10.075	1440	1443	1451	rVB6	5251	9859	0.56%	0.094%
49	10.151	1451	1456	1459	rBV3	2977	4718	0.27%	0.045%
50	10.269	1468	1476	1489	rBV	165379	316953	17.89%	3.025%
51	10.440	1501	1505	1507	rBV4	2667	3576	0.20%	0.034%
52	10.522	1513	1519	1525	rBV	40925	77565	4.38%	0.740%
53	10.616	1531	1535	1538	rVB3	5885	8966	0.51%	0.086%
54	10.669	1541	1544	1546	rVB2	2507	2281	0.13%	0.022%
55	10.704	1546	1550	1552	rBV3	7016	8633	0.49%	0.082%
56	10.746	1555	1557	1558	rVV	3048	2730	0.15%	0.026%
57	10.810	1560	1568	1574	rVV	449267	795542	44.90%	7.592%
58	10.875	1574	1579	1591	rVB	95875	170823	9.64%	1.630%
59	11.010	1599	1602	1604	rBV3	2715	3457	0.20%	0.033%
60	11.057	1606	1610	1619	rVV3	18216	37725	2.13%	0.360%
61	11.128	1621	1622	1624	rVV2	5121	3260	0.18%	0.031%
62	11.175	1624	1630	1637	rVV3	34660	71314	4.02%	0.681%
63	11.228	1637	1639	1644	rVV4	10253	17117	0.97%	0.163%
64	11.281	1644	1648	1655	rVV8	8082	17130	0.97%	0.163%
65	11.340	1655	1658	1661	rVV4	3030	4432	0.25%	0.042%
66	11.387	1661	1666	1676	rVB4	30929	67099	3.79%	0.640%
67	11.581	1692	1699	1709	rBV	356910	613955	34.65%	5.859%
68	11.716	1714	1722	1727	rBV7	18228	36101	2.04%	0.345%
69	11.769	1727	1731	1735	rBV5	12979	21503	1.21%	0.205%
70	11.816	1737	1739	1740	rBV2	7862	6591	0.37%	0.063%
71	11.928	1754	1758	1763	rBV3	5385	8865	0.50%	0.085%
72	11.987	1766	1768	1773	rVB3	3618	5009	0.28%	0.048%
73	12.093	1775	1786	1794	rBV4	35268	96762	5.46%	0.923%
74	12.257	1809	1814	1816	rBV5	6107	9907	0.56%	0.095%
75	12.298	1818	1821	1822	rBV3	8045	10369	0.59%	0.099%
76	12.487	1849	1853	1856	rBV5	6643	12731	0.72%	0.121%

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Instrument :
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 ClientSampleId :
 A46S5RE

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0 % of largest Peak

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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_D\Method\SFAMDLM061523SMA.M
 Title : SFAM01.0

77	12.516	1856	1858	1863	rVB5	7681	10845	0.61%	0.103%
78	12.575	1863	1868	1873	rVB5	14362	27019	1.52%	0.258%
79	12.651	1873	1881	1887	rBV	142282	261705	14.77%	2.497%
80	12.740	1891	1896	1901	rVB2	54434	96773	5.46%	0.924%
81	12.834	1904	1912	1914	rBV8	19981	50401	2.84%	0.481%
82	13.081	1951	1954	1958	rBV3	6173	7974	0.45%	0.076%
83	13.134	1958	1963	1965	rBV6	5592	10270	0.58%	0.098%
84	13.222	1972	1978	1979	rBV4	9932	16573	0.94%	0.158%
85	13.351	1994	2000	2003	rBV5	12971	27292	1.54%	0.260%
86	13.516	2021	2028	2036	rBV	358640	591936	33.41%	5.649%
87	13.822	2072	2080	2090	rVB2	356937	797025	44.98%	7.606%
88	14.051	2115	2119	2126	rVB4	10344	18034	1.02%	0.172%
89	14.216	2144	2147	2155	rVB7	5485	9612	0.54%	0.092%
90	14.345	2167	2169	2171	rBV3	5998	6707	0.38%	0.064%
91	14.616	2213	2215	2216	rBV2	3351	1575	0.09%	0.015%
92	14.634	2216	2218	2222	rVV4	2946	2406	0.14%	0.023%
93	14.775	2237	2242	2243	rBV5	17551	23045	1.30%	0.220%
94	15.098	2292	2297	2302	rBV5	15962	25405	1.43%	0.242%
95	15.486	2356	2363	2367	rVB3	10813	22492	1.27%	0.215%
96	15.522	2367	2369	2371	rBV2	3869	3242	0.18%	0.031%
97	15.781	2412	2413	2415	rBV2	2492	1770	0.10%	0.017%
98	16.045	2456	2458	2459	rBV2	3205	1791	0.10%	0.017%
99	16.281	2497	2498	2500	rBV2	3147	2065	0.12%	0.020%
100	16.775	2580	2582	2584	rBV3	2664	1957	0.11%	0.019%

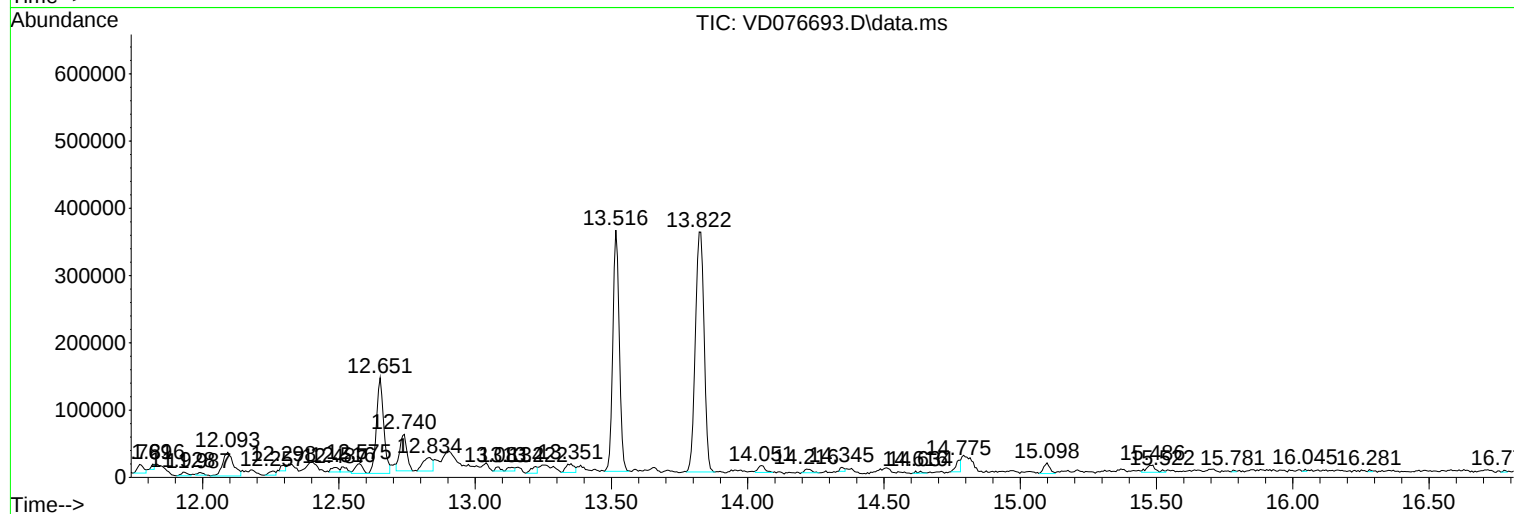
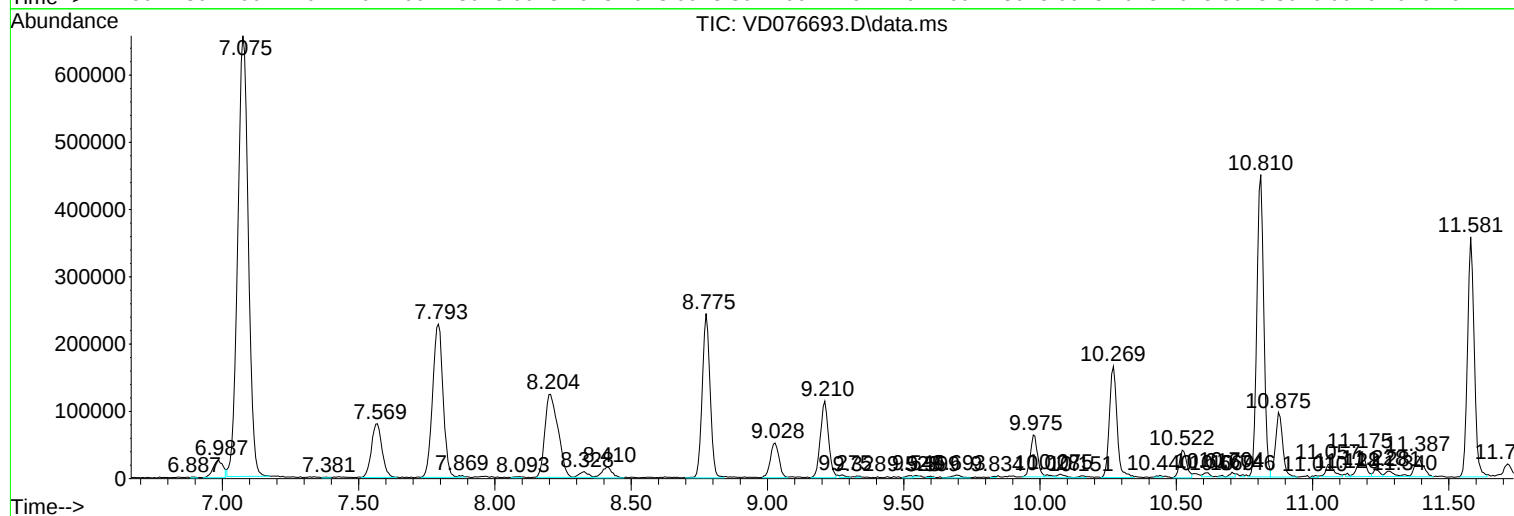
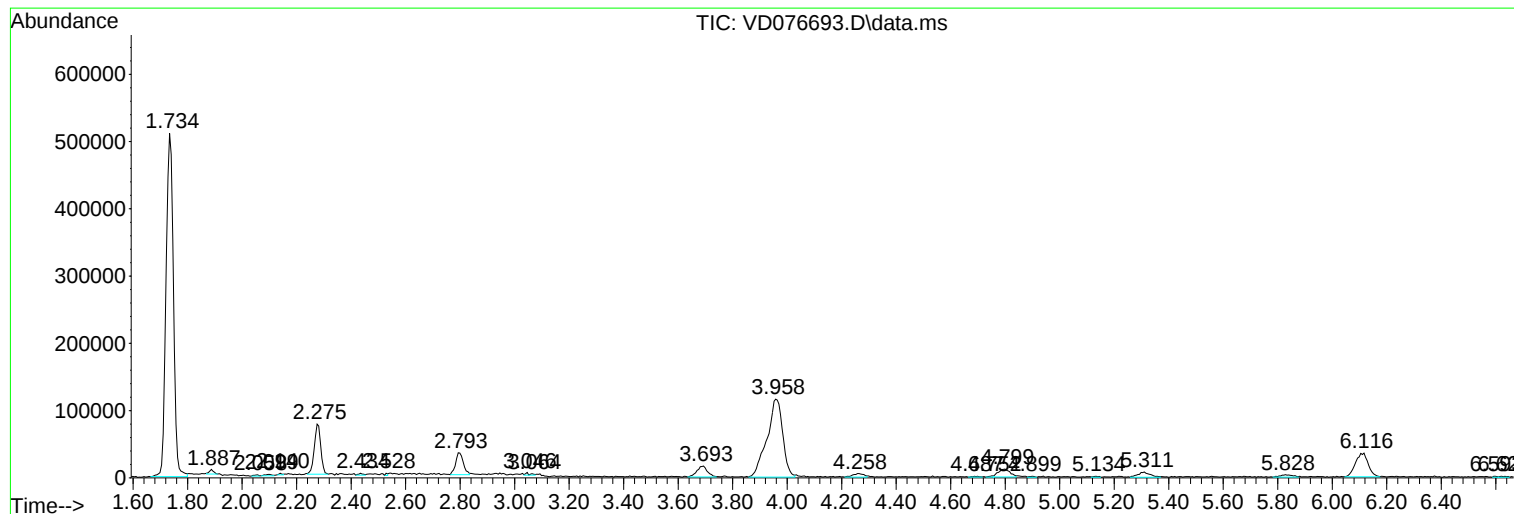
Sum of corrected areas: 10478776

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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
A46S5RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM061523SMA.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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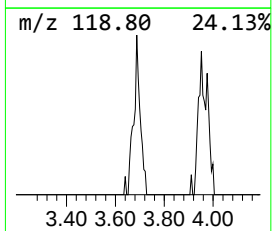
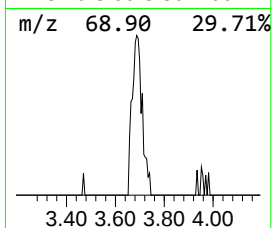
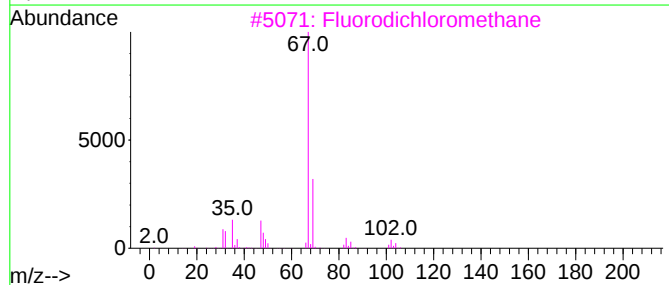
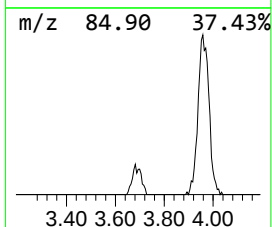
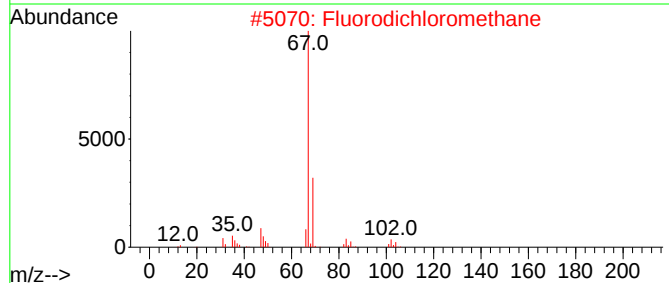
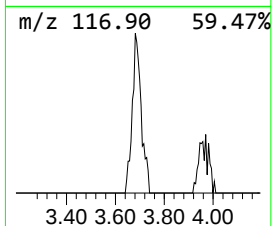
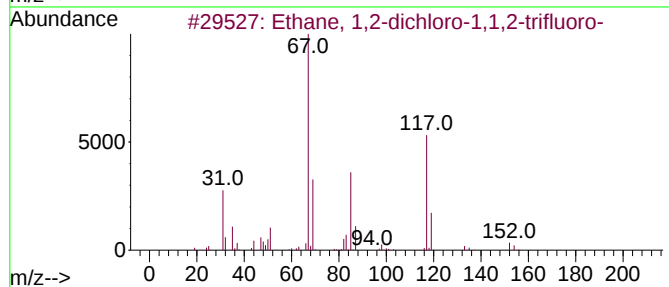
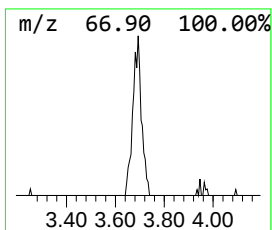
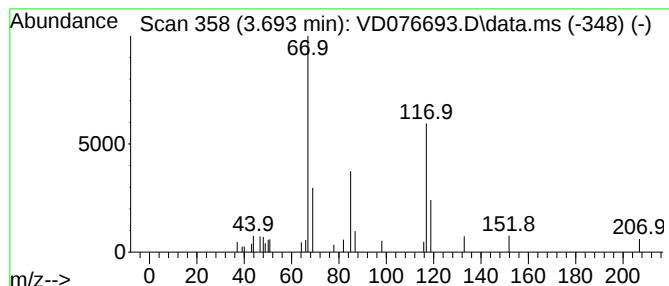
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM061523SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.693	2.17 ug/L	42788	1,4-Difluorobenzene	8.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethane, 1,2-dichloro-1,1,2-trifl...	152	C2HCl2F3	000354-23-4	38
2			Fluorodichloromethane	102	CHCl2F	000075-43-4	32
3			Fluorodichloromethane	102	CHCl2F	000075-43-4	23
4			Fluorodichloromethane	102	CHCl2F	000075-43-4	17
5			Propane, 1,3-dichloro-1,1,2,2,3-...	202	C3HCl2F5	000507-55-1	10



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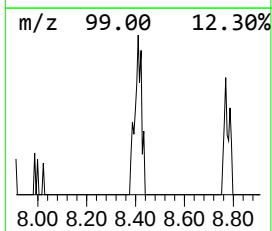
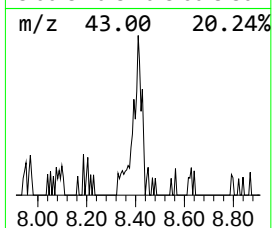
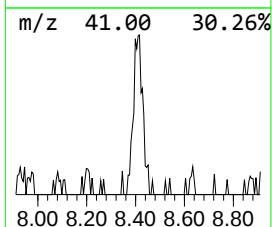
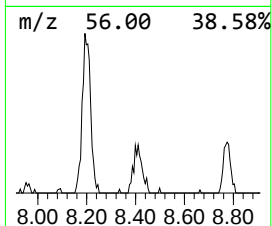
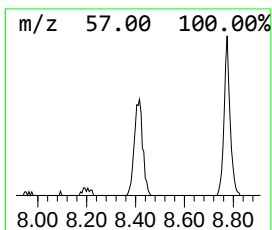
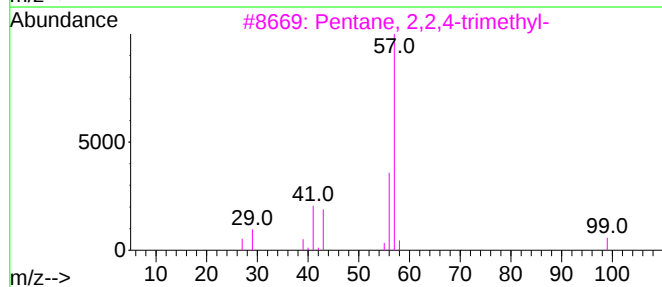
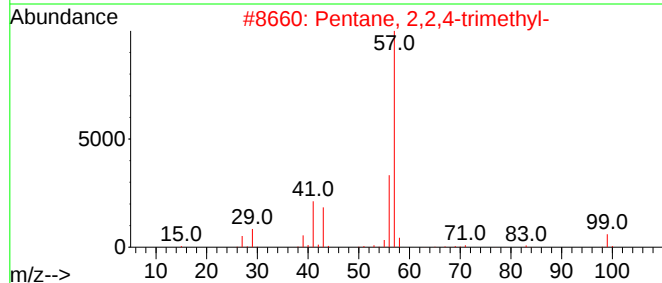
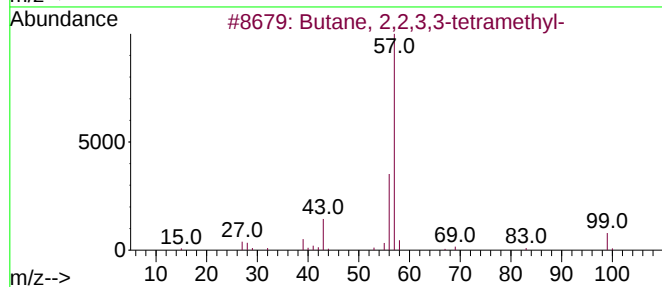
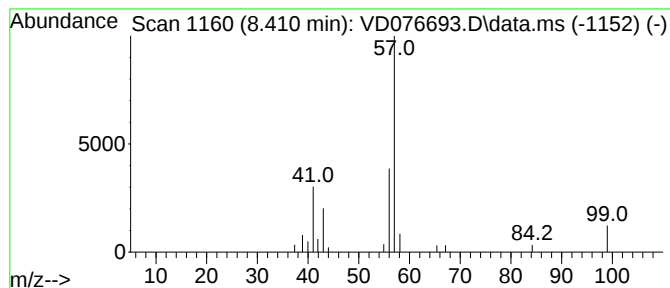
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM061523SMA.M
Quant Title : SFAM01.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.410	2.11 ug/L	41620	1,4-Difluorobenzene	8.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	56
2		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	56
3		Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	56
4		Butane, 2-azido-2,3,3-trimethyl-	141	C7H15N3	051677-41-9	50
5		Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	39



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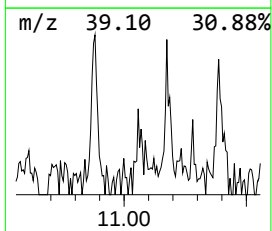
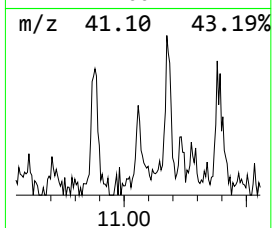
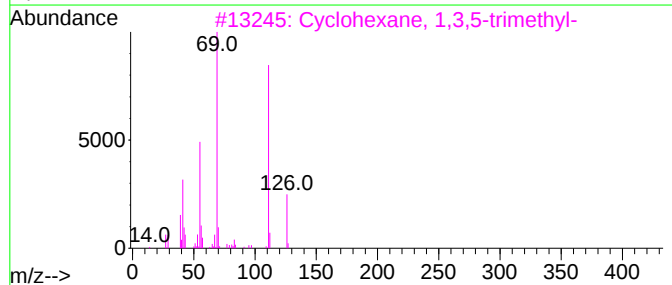
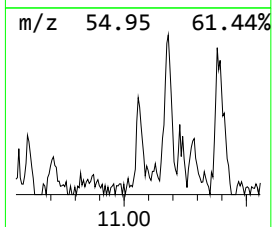
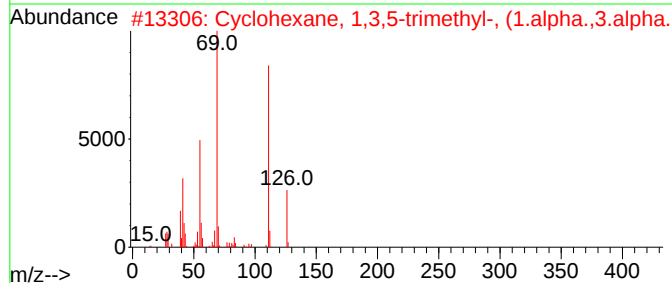
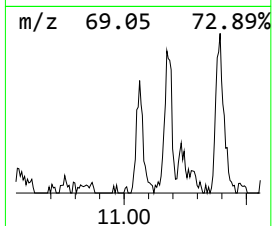
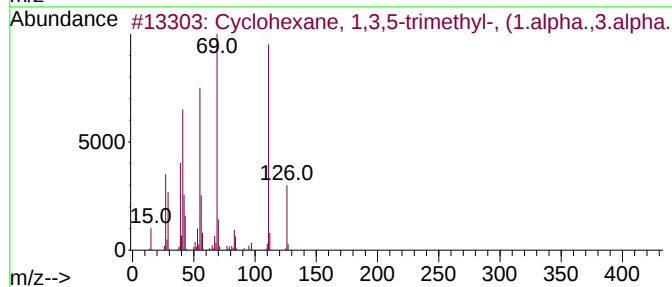
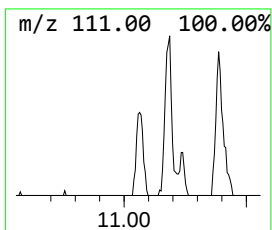
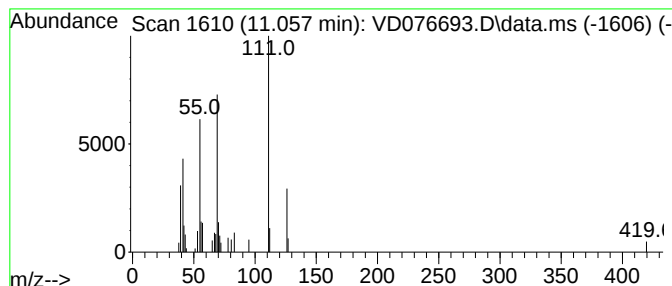
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Quant Title : SFAM01.0

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

Peak Number 6 (DEL) Alkane: Cyclic11.057 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.057	1.54 ug/L	37725	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	87
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	83
3			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	83
4			3-Buten-2-one, 3-methyl-, dimeth...	126	C7H14N2	075268-10-9	80
5			Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
A46S5RE

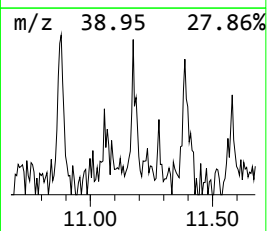
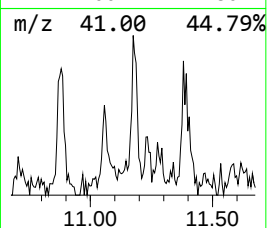
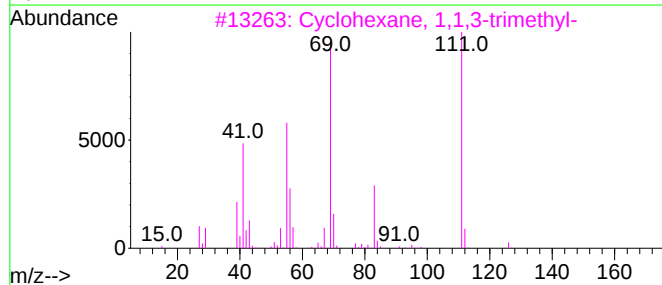
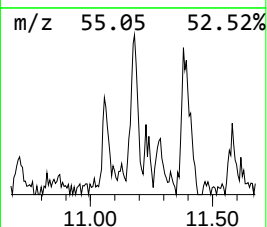
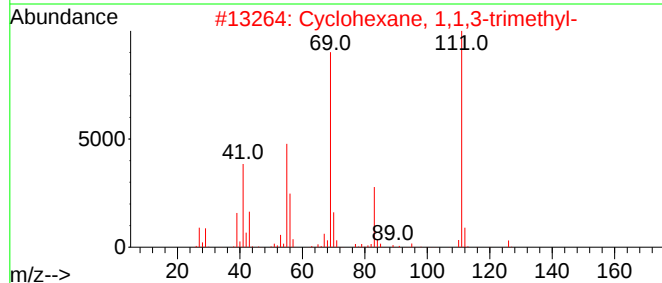
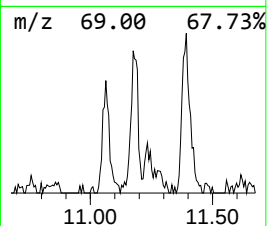
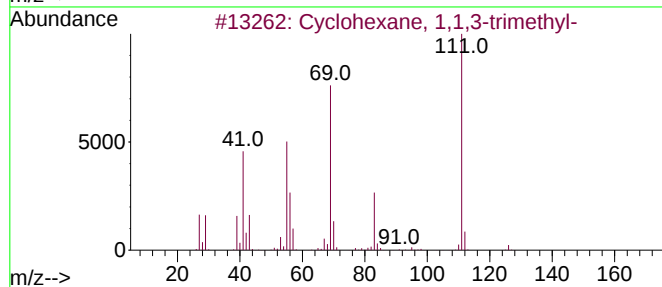
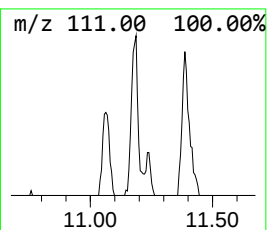
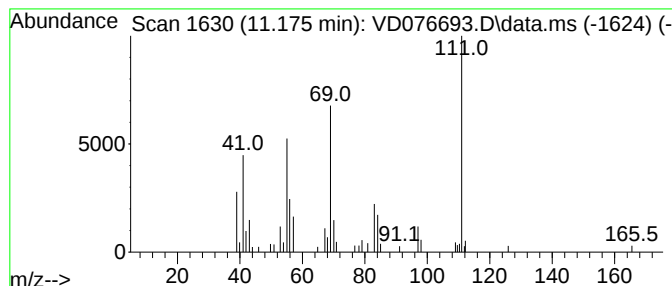
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Quant Title : SFAM01.0

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

Peak Number 7 (DEL) Alkane: Cyclic11.175 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.175	2.90 ug/L	71314	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	81
2			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	81
3			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	76
4			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	64
5			3'-Fluoroacetanilide	153	C8H8FNO	000351-28-0	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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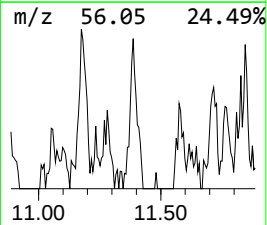
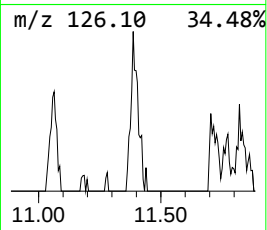
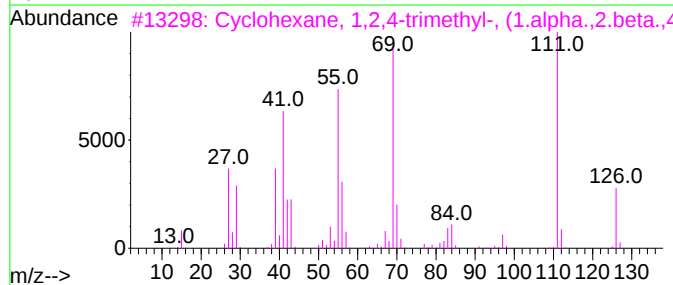
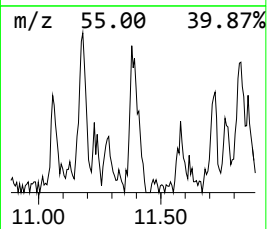
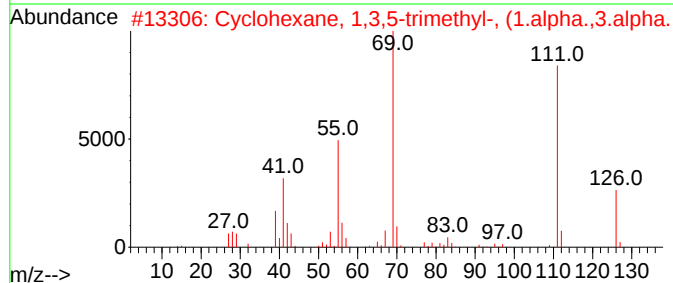
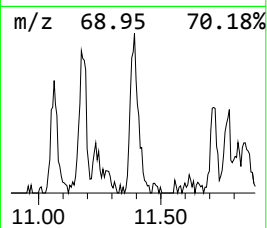
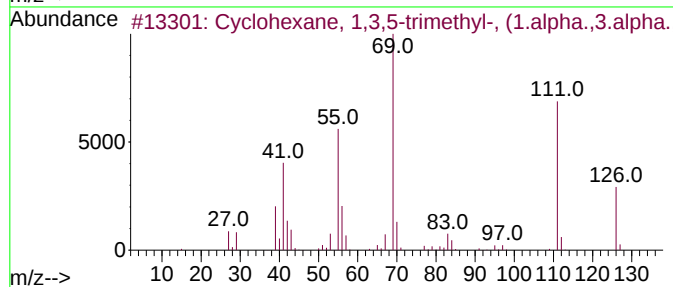
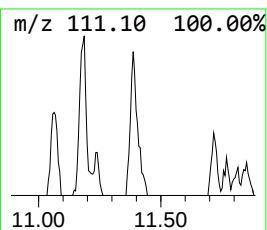
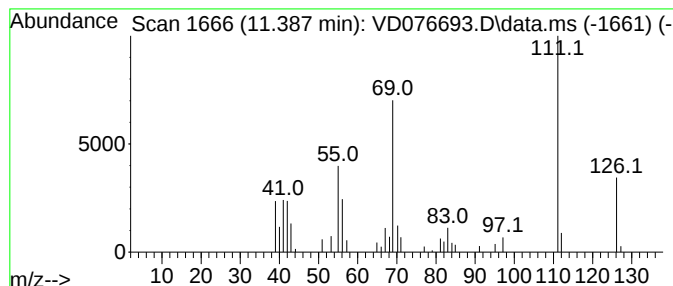
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Quant Title : SFAM01.0

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

Peak Number 8 (DEL) Alkane: Cyclic11.387 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.387	2.73 ug/L	67099	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	86
2			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	80
3			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	64
4			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	64
5			Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	59



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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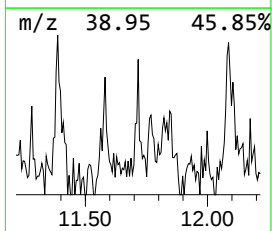
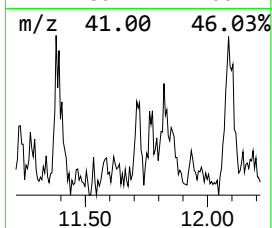
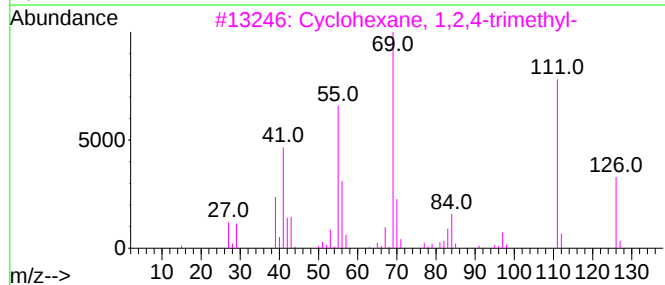
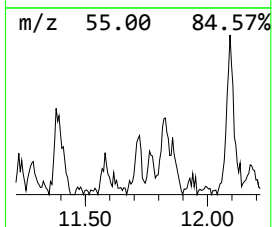
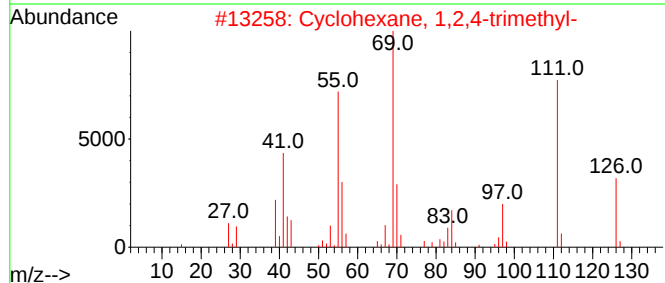
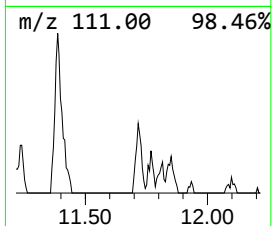
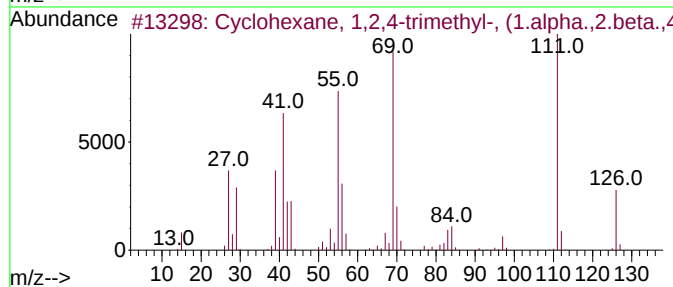
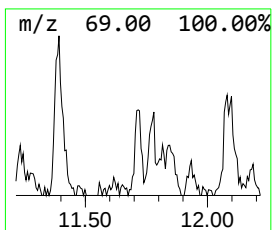
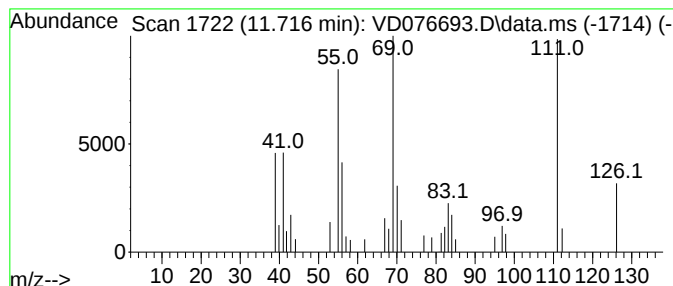
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Quant Title : SFAM01.0

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

Peak Number 9 (DEL) Alkane: Cyclic11.716 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.716	1.47 ug/L	36101	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	91
2			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
3			Cyclohexane, 1,2,4-trimethyl-	126	C9H18	002234-75-5	90
4			Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	81
5			Cyclohexane, 1,2,4-trimethyl-, (...	126	C9H18	007667-60-9	74



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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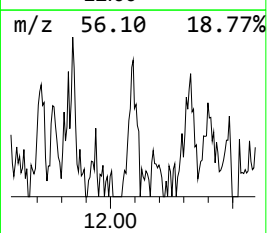
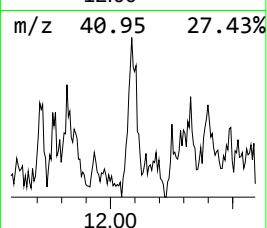
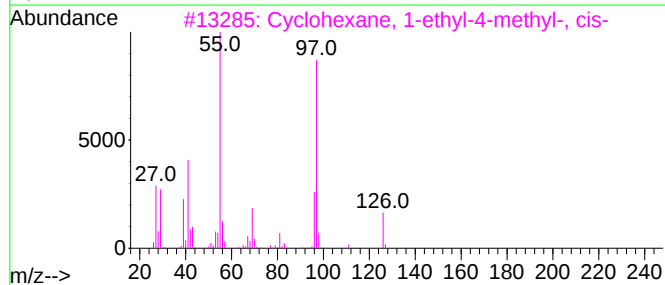
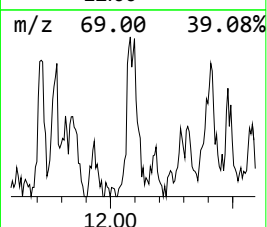
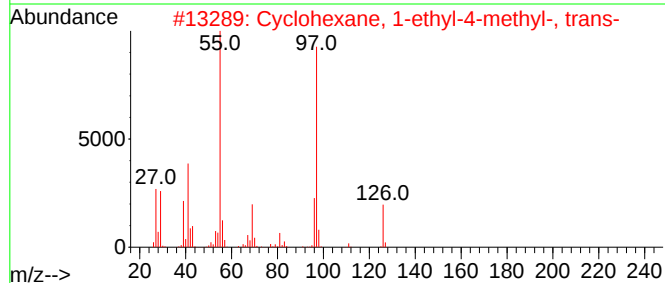
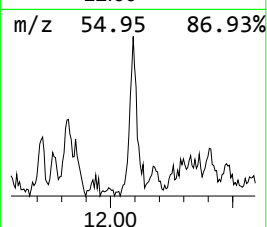
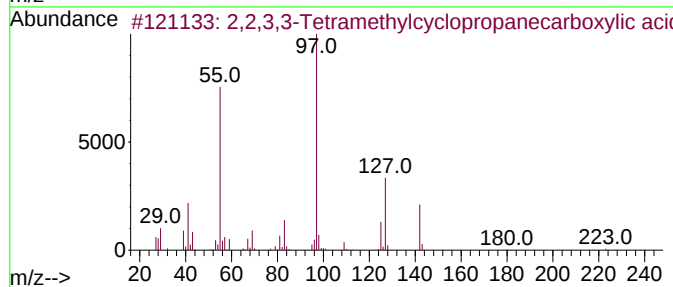
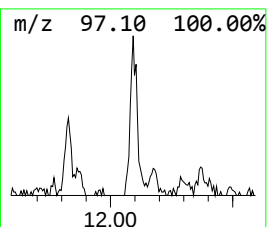
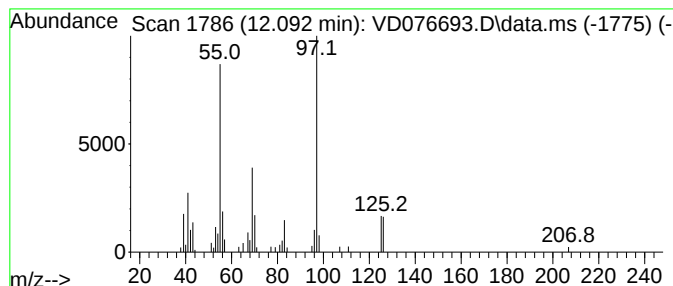
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Quant Title : SFAM01.0

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

Peak Number 10 2,2,3,3-Tetramethylcyclopro... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.092	3.94 ug/L	96762	Chlorobenzene-d5	11.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,3,3-Tetramethylcyclopropanec...	238	C15H26O2	1010221-97-5	78
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	72
3			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	72
4			Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	68
5			Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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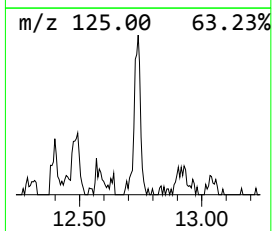
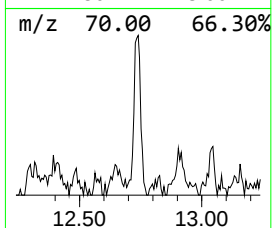
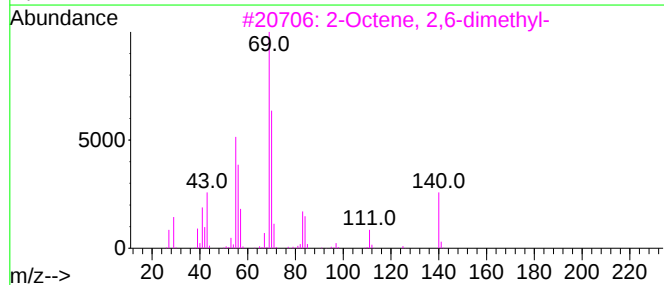
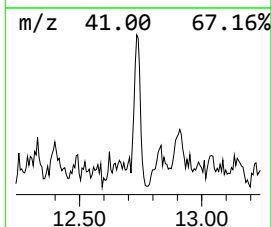
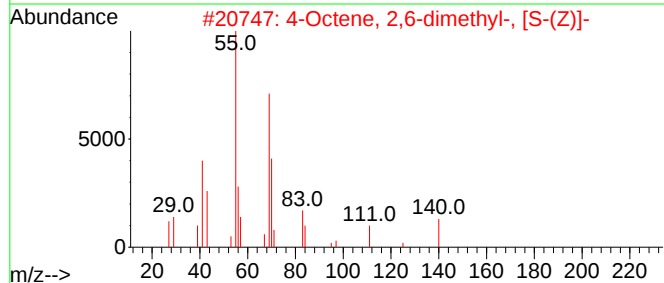
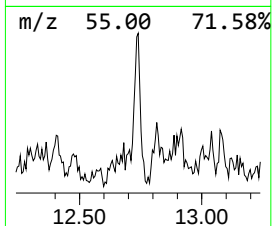
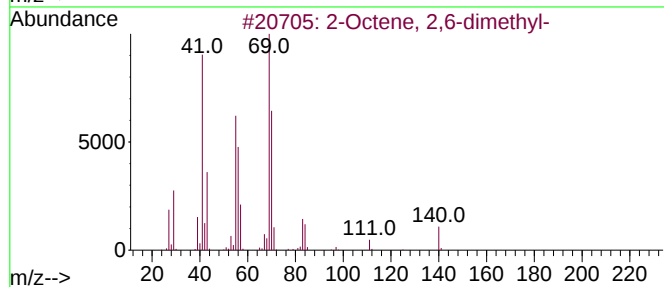
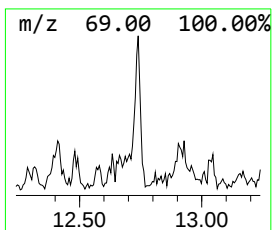
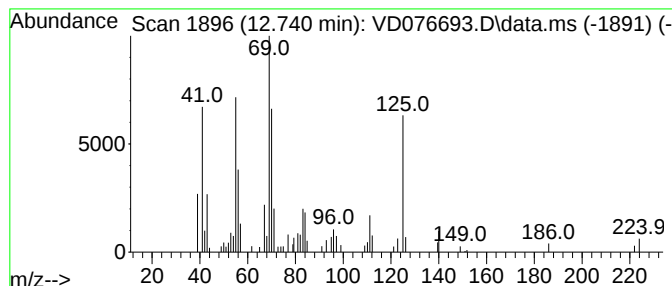
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Quant Title : SFAM01.0

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.740	4.09 ug/L	96773	1,4-Dichlorobenzene-d4	13.516

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	62
2		4-Octene, 2,6-dimethyl-, [S-(Z)]-	140	C10H20	062960-77-4	58
3		2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	53
4		3-Heptene, 2-methyl-	112	C8H16	017618-76-7	43
5		1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	43



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
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Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

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ClientSampleId :
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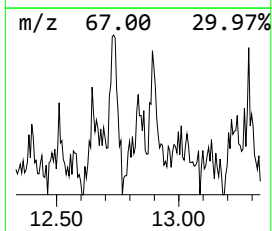
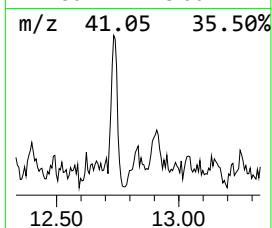
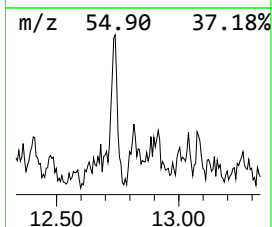
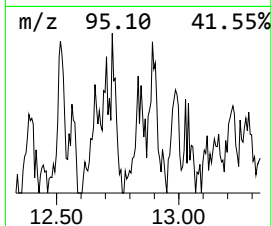
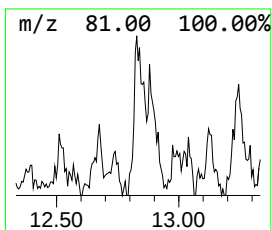
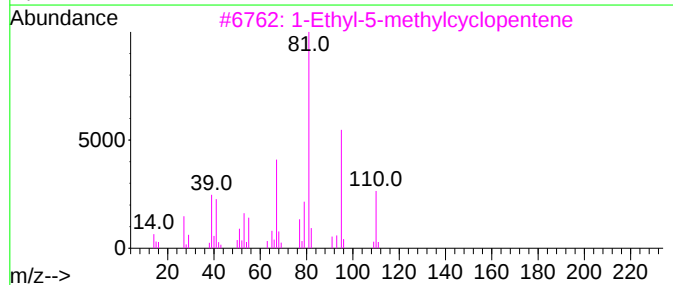
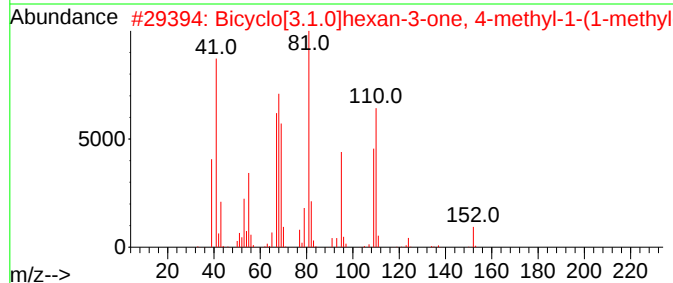
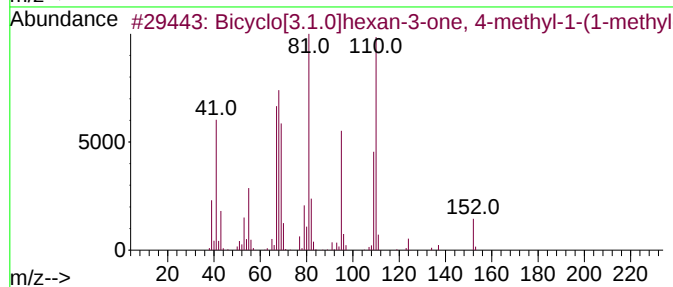
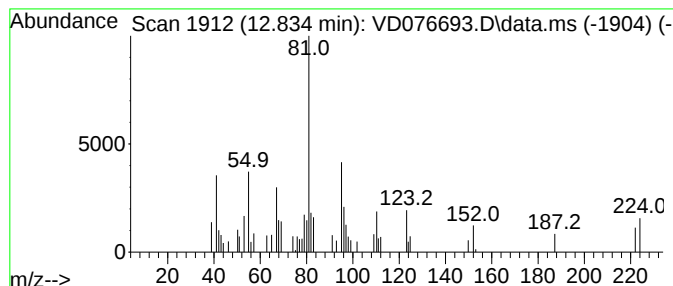
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Quant Title : SFAM01.0

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

Peak Number 12 Bicyclo[3.1.0]hexan-3-one, ... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.834	2.13 ug/L	50401	1,4-Dichlorobenzene-d4	13.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[3.1.0]hexan-3-one, 4-met...	152	C10H16O	000471-15-8	58
2			Bicyclo[3.1.0]hexan-3-one, 4-met...	152	C10H16O	001125-12-8	53
3			1-Ethyl-5-methylcyclopentene	110	C8H14	097797-57-4	53
4			Methyl ethyl cyclopentene	110	C8H14	019780-56-4	47
5			Cyclopentane, 1,3-dimethyl-2-(1-...	138	C10H18	061142-30-1	47



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD070323\
Data File : VD076693.D
Acq On : 03 Jul 2023 15:35
Operator : KP/SY
Sample : 03472-13RE
Misc : 5.90G/10.0ml/MSVOA_D/SOIL
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
A46S5RE

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM061523SMA.M
Quant Title : SFAM01.0

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	3.693	2.2	ug/L	42788	1	8.775	492395	25.0
(DEL) Alkane: S...	8.410	2.1	ug/L	41620	1	8.775	492395	25.0
(DEL) Alkane: C...	11.057	1.5	ug/L	37725	2	11.581	613955	25.0
(DEL) Alkane: C...	11.175	2.9	ug/L	71314	2	11.581	613955	25.0
(DEL) Alkane: C...	11.387	2.7	ug/L	67099	2	11.581	613955	25.0
(DEL) Alkane: C...	11.716	1.5	ug/L	36101	2	11.581	613955	25.0
2,2,3,3-Tetrame...	12.092	3.9	ug/L	96762	2	11.581	613955	25.0
(DEL) Alkane: S...	12.740	4.1	ug/L	96773	3	13.516	591936	25.0
Bicyclo[3.1.0]h...	12.834	2.1	ug/L	50401	3	13.516	591936	25.0