LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

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\SFAMWLM040722SMA.M
Title : SFAM01.0

Signal : TIC: VW022514.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.154	38	41	42	rBV3	2130	2682	0.01%	0.005%
2	2.355	69	74	82	rBV	149525	238153	0.83%	0.488%
3	2.886	156	161	172	rVB	82806	185241	0.65%	0.379%
4	3.392	242	244	247	rVB3	1169	1194	0.00%	0.002%
5	3.459	252	255	256	rBV3	819	659	0.00%	0.001%
6	3.581	273	275	276	rBV2	1263	812	0.00%	0.002%
7	3.605	276	279	283	rVB5	1380	1322	0.00%	0.003%
8	3.654	285	287	289	rBV2	1178	1131	0.00%	0.002%
9	3.733	296	300	301	rBV4	1237	1372	0.00%	0.003%
10	4.019	336	347	361	rBV3	149515	470016	1.64%	0.963%
11	4.221	378	380	382	rBV3	687	688	0.00%	0.001%
12	4.263	382	387	388	rBV5	1230	2181	0.01%	0.004%
13	4.361	399	403	406	rBV5	1628	2601	0.01%	0.005%
14	4.471	415	421	422	rVB7	1307	1721	0.01%	0.004%
15	4.532	429	431	433	rBV3	1269	1240	0.00%	0.003%
16	4.641	446	449	453	rBV5	1574	2455	0.01%	0.005%
17	4.672	453	454	455	rVV	1690	1010	0.00%	0.002%
18	4.702	458	459	462	rVV3	1645	1568	0.01%	0.003%
19	4.745	464	466	469	rVV3	1249	1610	0.01%	0.003%
20	4.769	469	470	473	rVV3	1657	1696	0.01%	0.003%
21	4.818	476	478	481	rVV3	1502	2242	0.01%	0.005%
22	4.916	485	494	508	rVV4	13708	52294	0.18%	0.107%
23	5.013	508	510	517	rVV6	1072	1679	0.01%	0.003%
24	5.105	524	525	527	rBV2	849	653	0.00%	0.001%
25	5.275	551	553	557	rVB5	718	812	0.00%	0.002%
26	5.361	560	567	570	rBV7	818	2145	0.01%	0.004%
27	5.385	570	571	574	rVB2	1070	879	0.00%	0.002%
28	5.409	574	575	576	rBV	900	645	0.00%	0.001%
29	5.525	593	594	596	rBV2	1086	903	0.00%	0.002%
30	5.574	600	602	604	rVB2	866	690	0.00%	0.001%
31	5.617	604	609	611	rBV6	1020	1497	0.01%	0.003%
32	5.720	624	626	632	rVB8	1076	2012	0.01%	0.004%
33	5.800	634	639	642	rBV4	533	872	0.00%	0.002%
34	5.879	649	652	653	rBV4	810	845	0.00%	0.002%
35	5.897	653	655	656	rBV2	1015	934	0.00%	0.002%
36	5.940	656	662	667	rBV8	3684	9755	0.03%	0.020%

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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

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\SFAMWLM040722SMA.M
Title : SFAM01.0

37	6.013	672	674	678	rVB3	1308	873	0.00%	0.002%
38	6.104	683	689	690	rBV5	1395	1824	0.01%	0.004%
39	6.141	693	695	697	rBV2	894	834	0.00%	0.002%
40	6.220	706	708	710	rVB3	848	727	0.00%	0.001%
41	6.293	718	720	721	rBV2	898	766	0.00%	0.002%
42	6.543	758	761	762	rVB3	726	647	0.00%	0.001%
43	6.562	762	764	767	rBV3	1194	1129	0.00%	0.002%
44	6.678	781	783	785	rVB2	1071	711	0.00%	0.001%
45	6.714	788	789	792	rVB3	877	804	0.00%	0.002%
46	6.763	792	797	800	rBV6	1249	1892	0.01%	0.004%
47	6.830	806	808	812	rVB5	820	891	0.00%	0.002%
48	6.873	812	815	817	rBV4	633	629	0.00%	0.001%
49	6.921	817	823	824	rVB4	911	1242	0.00%	0.003%
50	7.007	835	837	838	rBV2	694	667	0.00%	0.001%
51	7.080	840	849	863	rBV2	57894	167839	0.58%	0.344%
52	7.220	868	872	877	rBV7	847	1752	0.01%	0.004%
53	7.354	893	894	896	rBV2	981	754	0.00%	0.002%
54	7.647	931	942	958	rBV	284416	689933	2.40%	1.413%
55	7.757	958	960	961	rVV2	1065	855	0.00%	0.002%
56	7.775	961	963	968	rVB6	1273	1620	0.01%	0.003%
57	7.866	975	978	980	rBV3	1295	1920	0.01%	0.004%
58	8.159	1022	1026	1028	rBV5	1350	1882	0.01%	0.004%
59	8.183	1028	1030	1032	rBV3	823	1007	0.00%	0.002%
60	8.275	1034	1045	1062	rBV2	496002	1536493	5.35%	3.147%
61	8.842	1129	1138	1148	rBV	458227	921757	3.21%	1.888%
62	9.037	1164	1170	1171	rBV5	1156	2173	0.01%	0.004%
63	9.073	1174	1176	1177	rBV2	1803	1263	0.00%	0.003%
64	9.140	1185	1187	1190	rBV4	1202	1461	0.01%	0.003%
65	9.275	1198	1209	1218	rBV	355959	734623	2.56%	1.505%
66	10.031	1328	1333	1343	rVB	215819	402099	1.40%	0.824%
67	10.323	1374	1381	1388	rBV	723766	1273150	4.43%	2.608%
68	10.384	1389	1391	1398	rVB2	12608	20568	0.07%	0.042%
69	10.579	1407	1423	1428	rBV	153283	314805	1.10%	0.645%
70	11.628	1588	1595	1605	rVB	697685	1190614	4.15%	2.439%
71	11.738	1609	1613	1616	rVB6	3377	5863	0.02%	0.012%
72	12.061	1663	1666	1669	rBV5	1181	2114	0.01%	0.004%
73	12.207	1689	1690	1691	rBV	1430	996	0.00%	0.002%
74	12.689	1762	1769	1775	rBV	357071	594135	2.07%	1.217%
75	12.945	1806	1811	1812	rBV4	8074	11193	0.04%	0.023%
76	13.256	1845	1862	1880	rVB2	5951338	28711176	100.00%	58.809%

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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

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

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Peak separation: 5

Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Title : SFAM01.0

77	13.554	1904	1911	1918	rBV	854955	1360984	4.74%	2.788%
78	13.847	1952	1959	1967	rBV	784791	1337161	4.66%	2.739%
79	14.036	1986	1990	1992	rBV3	16625	24869	0.09%	0.051%
80	14.066	1992	1995	1999	rVB	24484	36131	0.13%	0.074%
81	14.213	2014	2019	2024	rBV2	19379	31488	0.11%	0.064%
82	14.414	2049	2052	2054	rBV4	6889	9154	0.03%	0.019%
83	14.463	2055	2060	2072	rVV	149472	256851	0.89%	0.526%
84	14.579	2075	2079	2083	rVB3	9286	13950	0.05%	0.029%
85	14.731	2096	2104	2113	rBV3	803635	1306767	4.55%	2.677%
86	14.816	2114	2118	2126	rVB2	33787	58744	0.20%	0.120%
87	15.005	2142	2149	2159	rBV2	51025	85363	0.30%	0.175%
88	15.176	2175	2177	2178	rBV2	2438	2260	0.01%	0.005%
89	15.219	2178	2184	2193	rVV9	15771	37924	0.13%	0.078%
90	15.334	2193	2203	2212	rVB5	32580	100450	0.35%	0.206%
91	15.548	2221	2238	2244	rBV	843018	1564440	5.45%	3.204%
92	15.603	2244	2247	2249	rVV3	11327	16594	0.06%	0.034%
93	15.725	2262	2267	2273	rVV7	8806	17529	0.06%	0.036%
94	15.786	2275	2277	2281	rVB5	2355	2540	0.01%	0.005%
95	15.859	2282	2289	2297	rVB	38546	78802	0.27%	0.161%
96	16.097	2325	2328	2335	rVB9	5621	11759	0.04%	0.024%
97	16.261	2349	2355	2362	rVB10	9792	22591	0.08%	0.046%
98	16.493	2378	2393	2400	rBV2	432681	1132418	3.94%	2.320%
99	16.609	2407	2412	2420	rVB2	59615	128140	0.45%	0.262%
100	16.724	2421	2431	2441	rBV	1622099	3580250	12.47%	7.333%

Sum of corrected areas: 48821049

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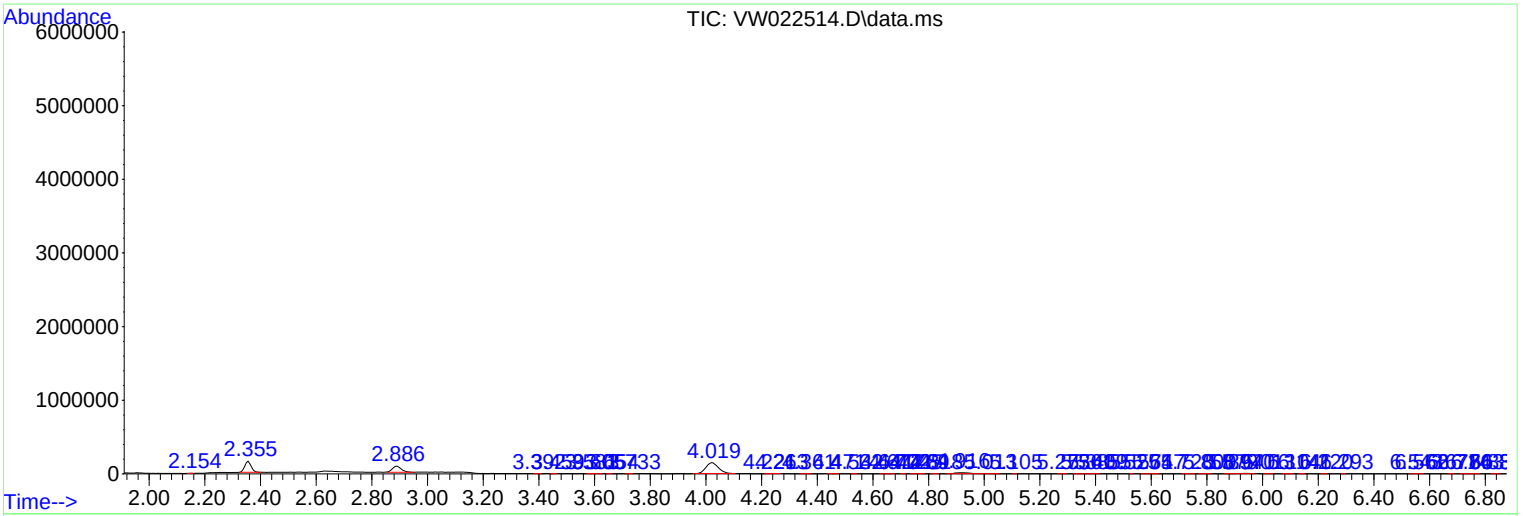
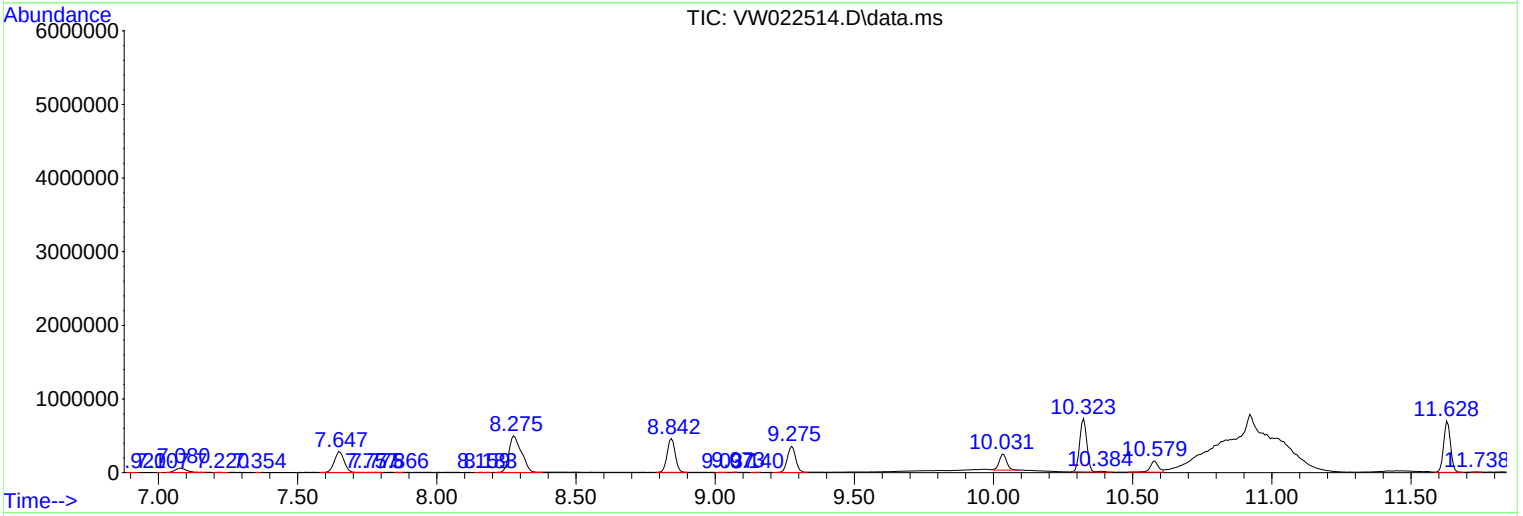
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

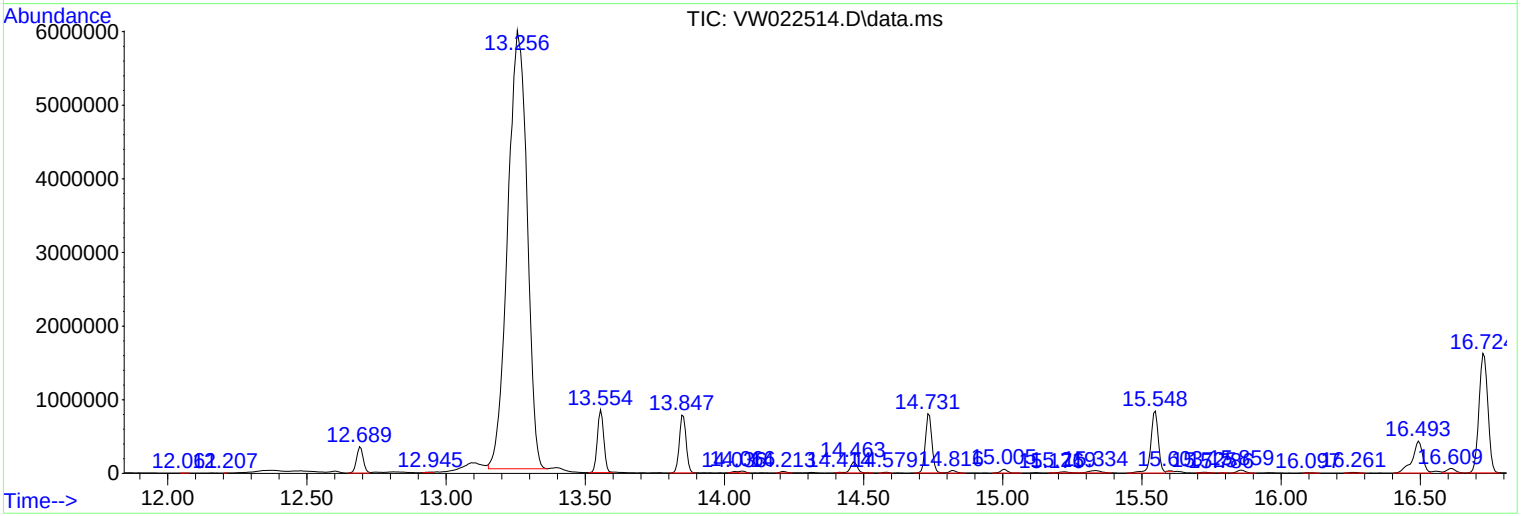
Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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







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Data File : VW022514.D
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Operator : SY/VA
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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

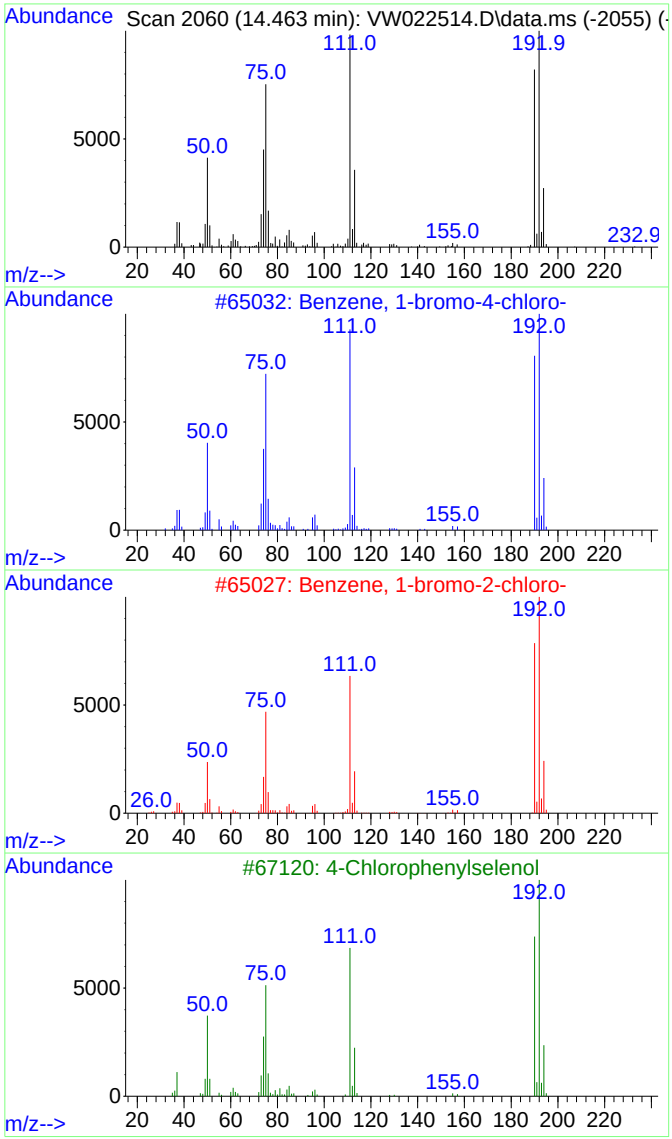
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

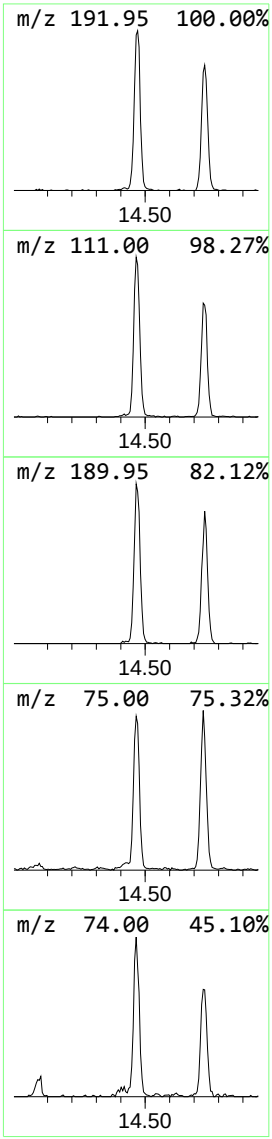
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-bromo-4-chloro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	4.72 ug/L	256851	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-bromo-4-chloro-	190	C6H4BrCl	000106-39-8	99
2	Benzene, 1-bromo-2-chloro-	190	C6H4BrCl	000694-80-4	98
3	4-Chlorophenylselenol	192	C6H5ClSe	016645-10-6	98
4	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	97
5	Benzene, 1-bromo-3-chloro-	190	C6H4BrCl	000108-37-2	96





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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

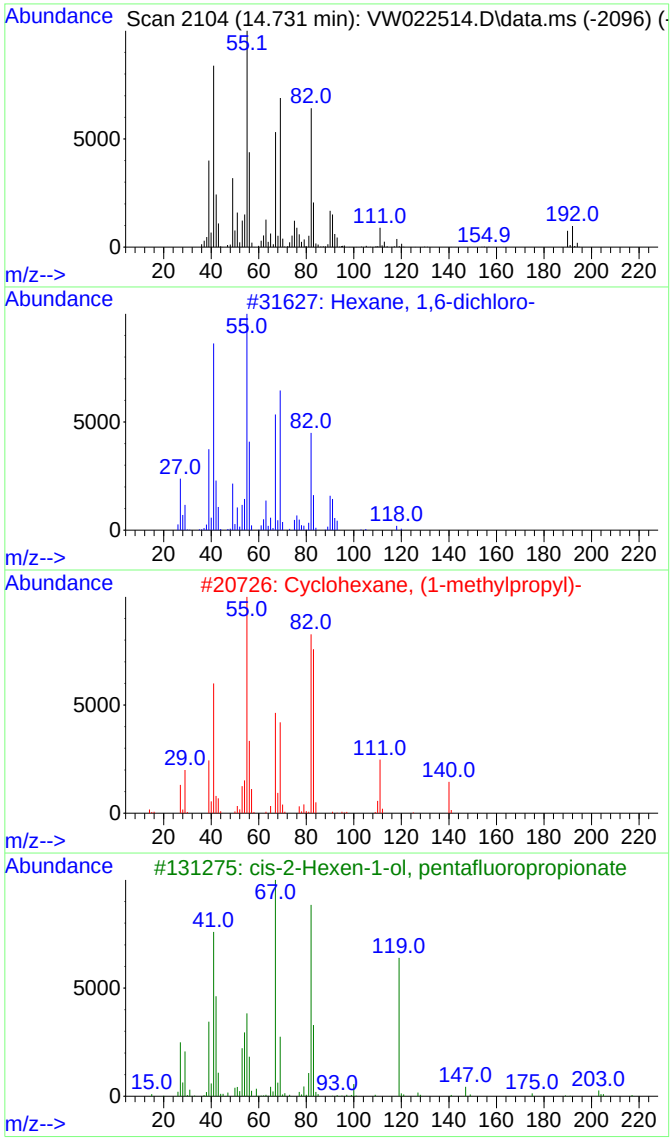
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

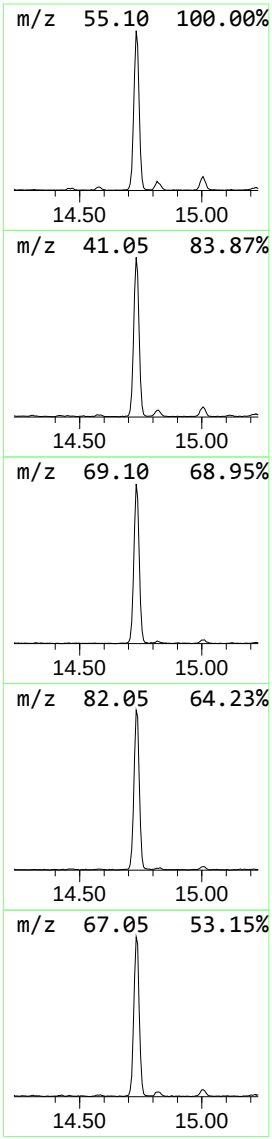
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Hexane, 1,6-dichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.731	24.00 ug/L	1306770	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1,6-dichloro-	154	C6H12Cl2	002163-00-0	90
2	Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	38
3	cis-2-Hexen-1-ol, pentafluoropro...	246	C9H11F5O2	1000352-74-1	35
4	Fumaric acid, cis-hex-3-enyl oct...	450	C28H50O4	1000348-87-3	35
5	Fumaric acid, hexadecyl trans-he...	422	C26H46O4	1000348-89-6	35





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Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

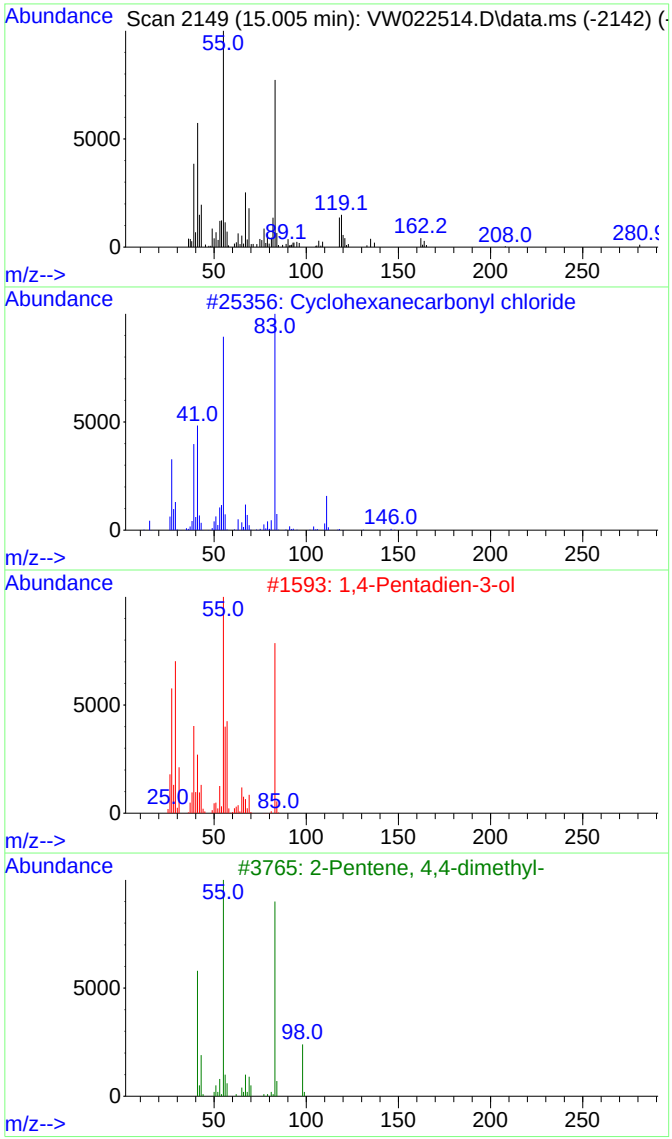
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

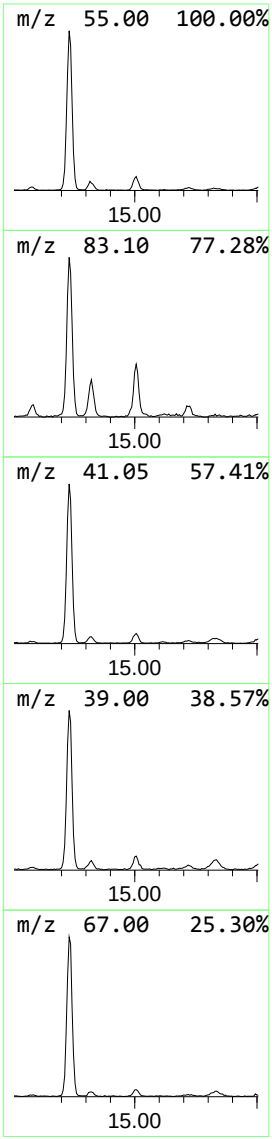
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclohexanecarbonyl chloride Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.005	1.57 ug/L	85363	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanecarbonyl chloride	146	C7H11ClO	002719-27-9	59
2	1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	58
3	2-Pentene, 4,4-dimethyl-	98	C7H14	026232-98-4	53
4	1-Hexene, 6-bromo-	162	C6H11Br	002695-47-8	53
5	Cyclopentane acetic acid hydrazide	142	C7H14N2O	020287-25-6	53





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Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

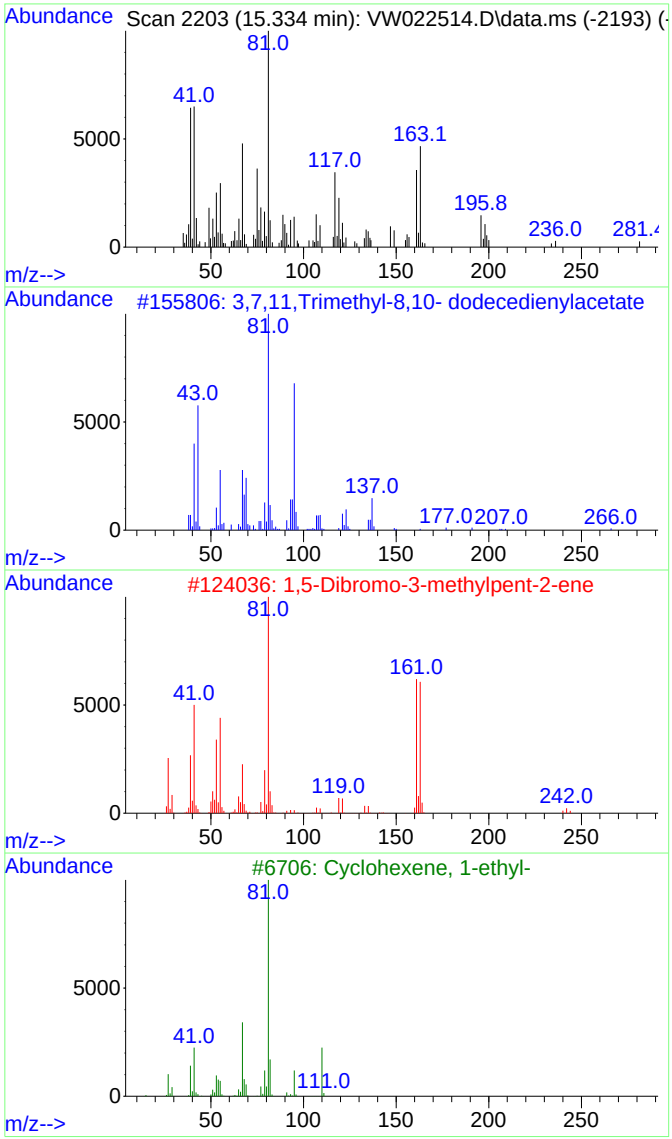
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

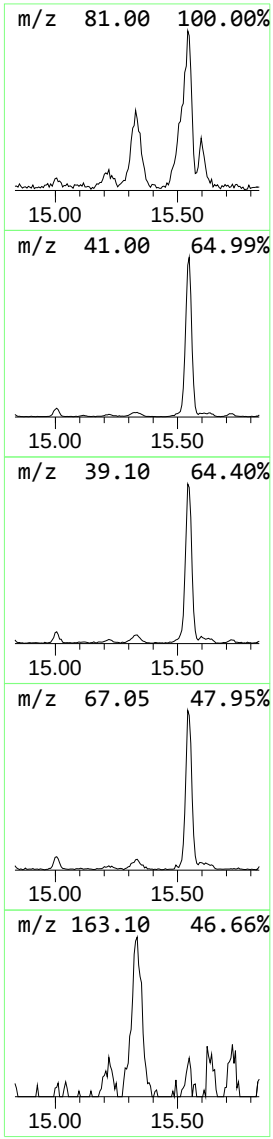
****

Peak Number 6 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.335	1.85 ug/L	100450	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,7,11,Trimethyl-8,10- dodecedie...	266	C17H30O2	1000374-10-3	35
2	1,5-Dibromo-3-methylpent-2-ene	240	C6H10Br2	133545-67-2	32
3	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	27
4	Cyclohexene, 3-methyl-	96	C7H12	000591-48-0	25
5	1-Heptyne	96	C7H12	000628-71-7	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
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Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

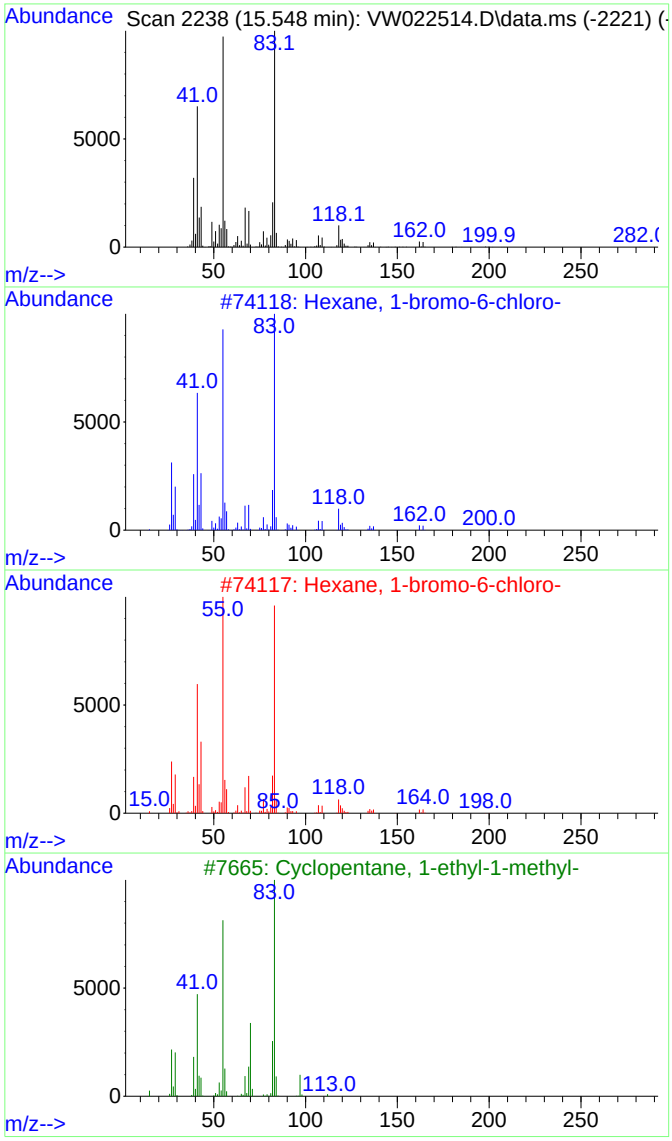
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

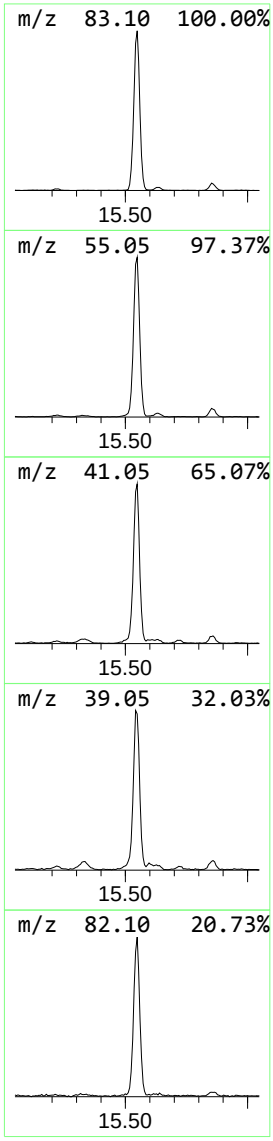
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Hexane, 1-bromo-6-chloro- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.548	28.74 ug/L	1564440	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	95
2	Hexane, 1-bromo-6-chloro-	198	C6H12BrCl	006294-17-3	72
3	Cyclopentane, 1-ethyl-1-methyl-	112	C8H16	016747-50-5	59
4	Pentanol, 4-methyl-4-nitro-	147	C6H13NO3	005215-92-9	50
5	Cyclohexanecarboxylic acid, ethe...	154	C9H14O2	004840-76-0	50





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Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

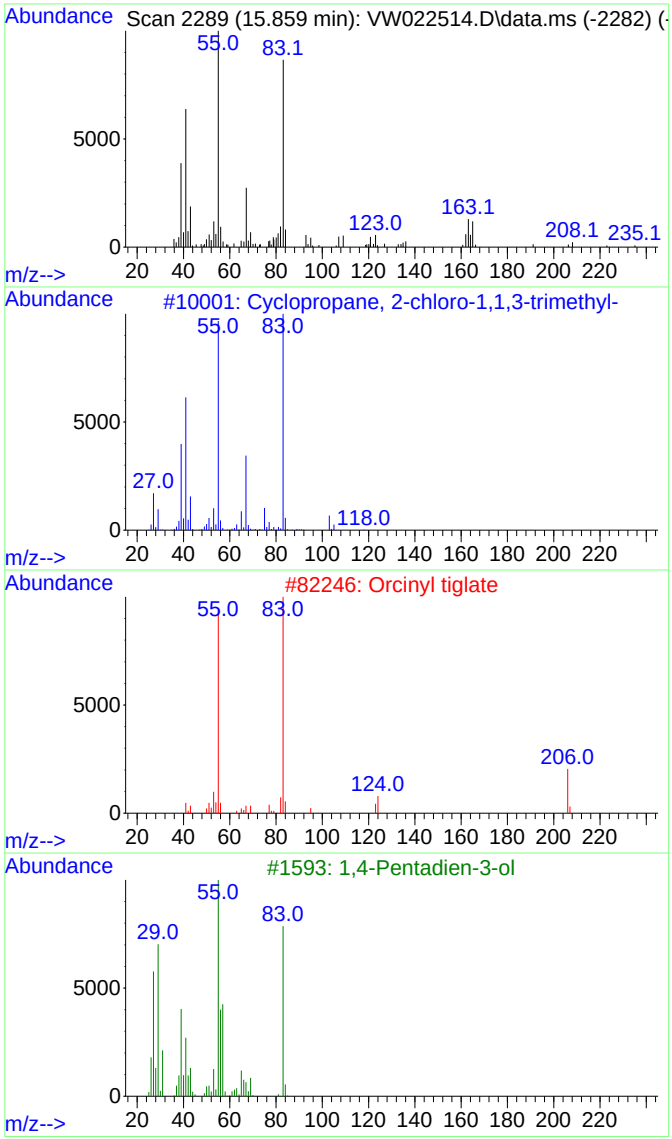
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

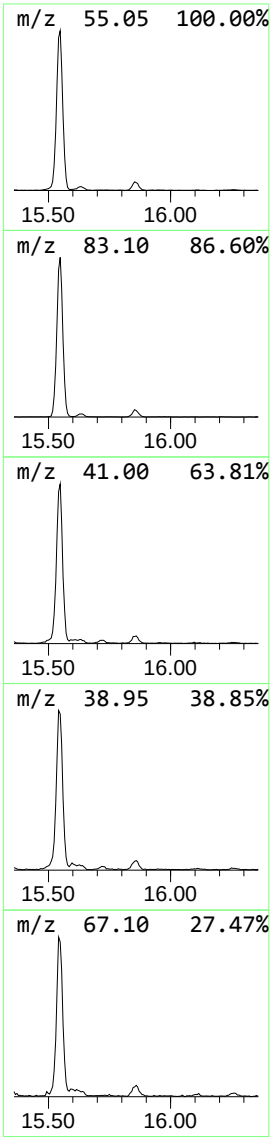
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Cyclopropane, 2-chloro-1,1,... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.859	1.45 ug/L	78802	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopropane, 2-chloro-1,1,3-tri...	118	C6H11Cl	098485-99-5	64
2	Orcinyl tiglate	206	C12H14O3	1000383-59-3	62
3	1,4-Pentadien-3-ol	84	C5H8O	000922-65-6	58
4	Octadienyl angelate, 2E, 4E-	208	C13H20O2	1000383-60-4	53
5	Hexane, 1,5-dibromo-	242	C6H12Br2	000627-96-3	52





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

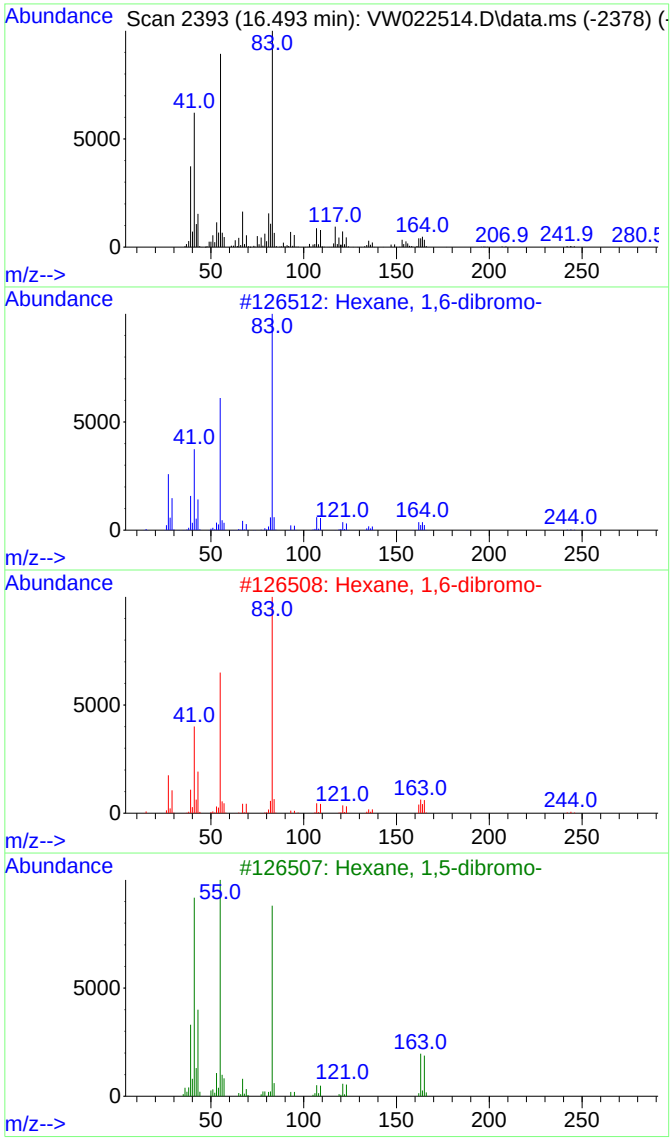
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

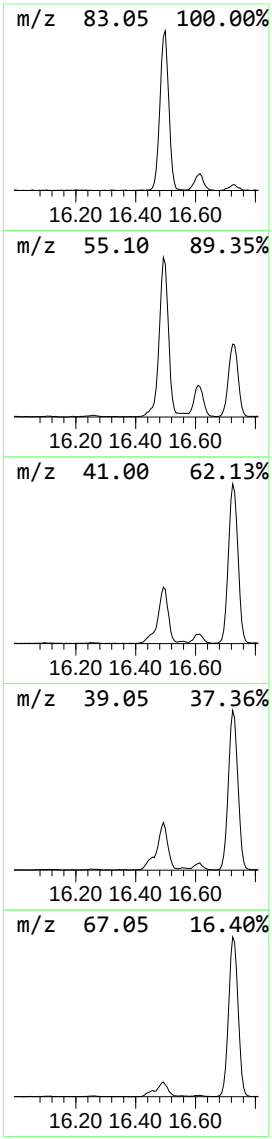
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Hexane, 1,6-dibromo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.493	20.80 ug/L	1132420	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	95
2	Hexane, 1,6-dibromo-	242	C6H12Br2	000629-03-8	91
3	Hexane, 1,5-dibromo-	242	C6H12Br2	000627-96-3	64
4	1-Hexene, 6-bromo-	162	C6H11Br	002695-47-8	59
5	Cumenyl angelate, o-	218	C14H18O2	1000383-67-2	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

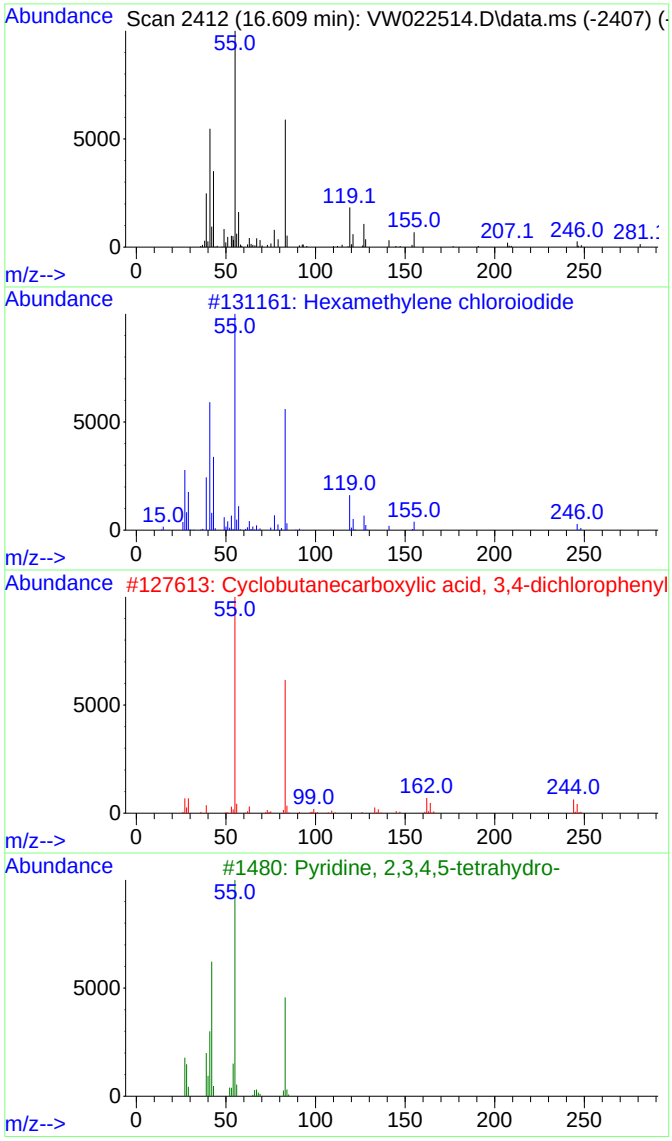
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

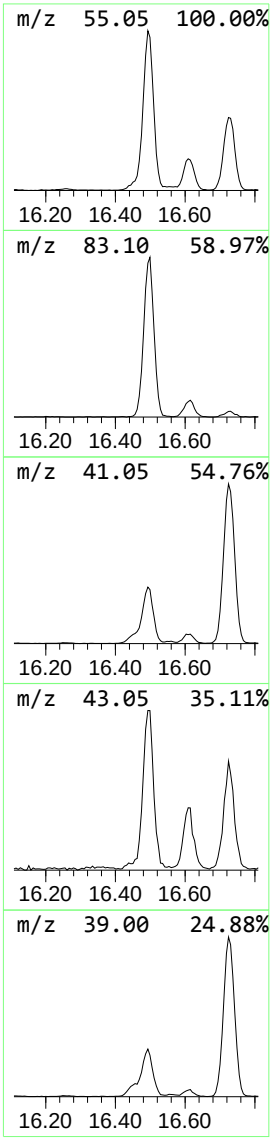
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Hexamethylene chloroiodide Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.609	2.35 ug/L	128140	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexamethylene chloroiodide	246	C6H12ClI	034683-73-3	97
2	Cyclobutanecarboxylic acid, 3,4-...	244	C11H10Cl2O2	1000307-44-1	47
3	Pyridine, 2,3,4,5-tetrahydro-	83	C5H9N	000505-18-0	43
4	Oxalic acid, cyclohexyl hexyl ester	256	C14H24O4	1000309-30-8	43
5	Cyclohexane, nitro-	129	C6H11NO2	001122-60-7	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

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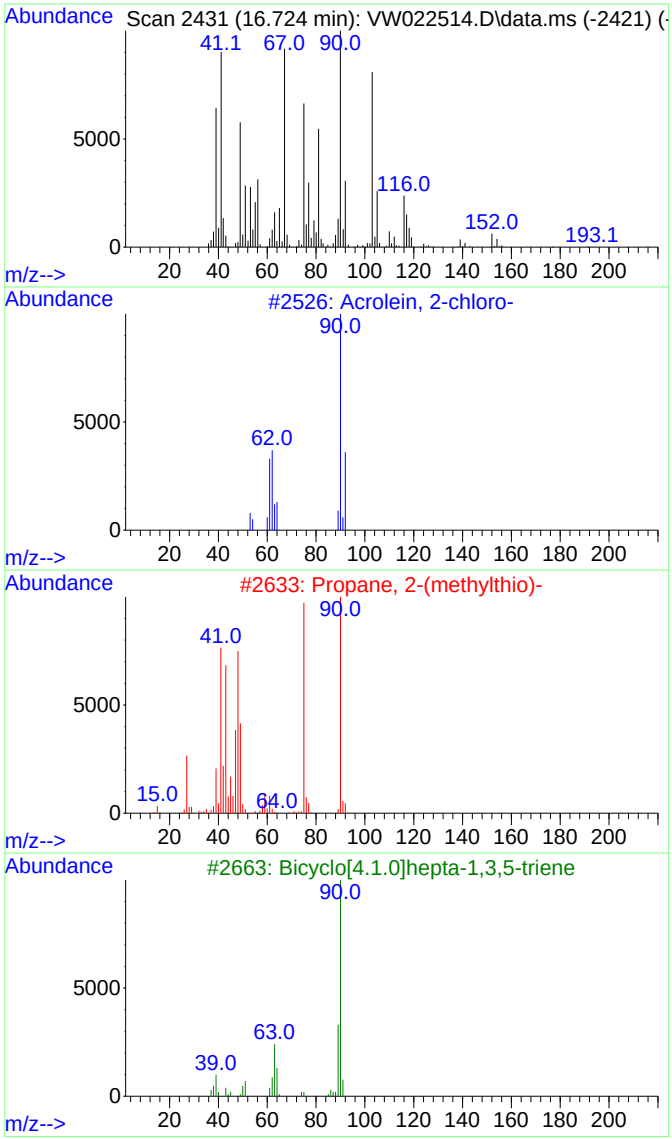
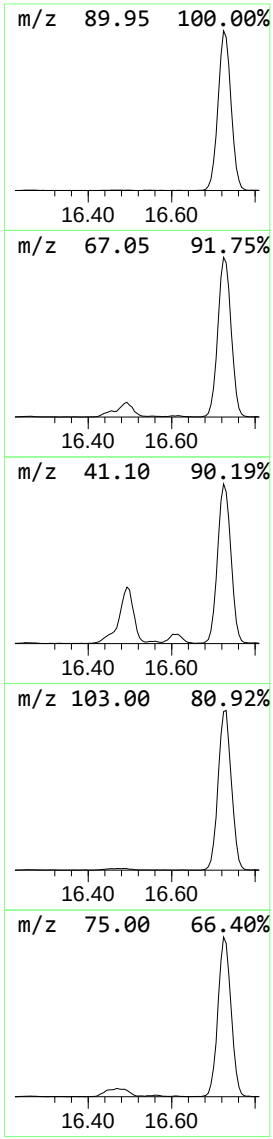
Peak Number 11 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.724	65.77 ug/L	3580250	1,4-Dichlorobenzene-d4	13.554

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acrolein, 2-chloro-	90	C3H3ClO	000683-51-2	43
2	Propane, 2-(methylthio)-	90	C4H10S	001551-21-9	32
3	Bicyclo[4.1.0]hepta-1,3,5-triene	90	C7H6	004646-69-9	30
4	2,2-Bis(chloromethyl)-1-propanol	156	C5H10Cl2O	005355-54-4	25
5	1H-Benzotriazole, 1-chloro-	153	C6H4ClN3	021050-95-3	12





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_W\Data\VW041222\
Data File : VW022514.D
Acq On : 12 Apr 2022 17:58
Operator : SY/VA
Sample : VIBLK423
Misc : 5.00g/10mL/MSVOA_W/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_W
ClientSampleId :
VIBLK423

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_W\Method\SFAMWLM040722SMA.M
Quant Title : SFAM01.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Benzene, 1-brom...	14.463	4.7	ug/L	256851	3	13.554	1360980	25.0
Hexane, 1,6-dic...	14.731	24.0	ug/L	1306770	3	13.554	1360980	25.0
Cyclohexanecarb...	15.005	1.6	ug/L	85363	3	13.554	1360980	25.0
unknown-01	15.335	1.9	ug/L	100450	3	13.554	1360980	25.0
Hexane, 1-bromo...	15.548	28.7	ug/L	1564440	3	13.554	1360980	25.0
Cyclopropane, 2...	15.859	1.5	ug/L	78802	3	13.554	1360980	25.0
Hexane, 1,6-dib...	16.493	20.8	ug/L	1132420	3	13.554	1360980	25.0
Hexamethylene c...	16.609	2.4	ug/L	128140	3	13.554	1360980	25.0
unknown-02	16.724	65.8	ug/L	3580250	3	13.554	1360980	25.0