

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
 Data File : VD077532.D
 Acq On : 09 Nov 2023 13:30
 Operator : JC/SY
 Sample : 05296-01
 Misc : 5.95G/10ml/MSVOA_D/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 BH3R1

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\SFAMDLM110923SMA.M
 Title : SFAM01.0

Signal : TIC: VD077532.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.746	19	27	43	rVB	2778923	4330121	3.17%	0.634%
2	1.893	49	52	60	rVB4	7434	14038	0.01%	0.002%
3	2.063	78	81	84	rBV2	2362	3793	0.00%	0.001%
4	2.152	92	96	99	rBV2	4276	7226	0.01%	0.001%
5	2.275	111	117	126	rBV	257975	394914	0.29%	0.058%
6	2.422	137	142	151	rVB	15847	28545	0.02%	0.004%
7	2.805	200	207	218	rBV	193511	358799	0.26%	0.053%
8	3.275	282	287	292	rBV5	1983	4613	0.00%	0.001%
9	3.922	386	397	408	rVV	244633	659062	0.48%	0.096%
10	4.028	408	415	422	rVV2	22807	63202	0.05%	0.009%
11	4.810	534	548	559	rBV3	26102	93144	0.07%	0.014%
12	5.316	628	634	645	rVV4	7727	22552	0.02%	0.003%
13	5.681	690	696	701	rBV5	2482	5411	0.00%	0.001%
14	5.840	715	723	734	rBV4	5409	18016	0.01%	0.003%
15	6.993	909	919	923	rBV2	42791	113924	0.08%	0.017%
16	7.081	923	934	968	rVB	12621902	33049895	24.17%	4.838%
17	7.569	1007	1017	1031	rVB	267993	684886	0.50%	0.100%
18	8.204	1115	1125	1142	rBV2	586284	1806911	1.32%	0.264%
19	8.640	1193	1199	1206	rVV8	4873	10716	0.01%	0.002%
20	8.781	1213	1223	1233	rBV	629349	1302918	0.95%	0.191%
21	9.028	1250	1265	1287	rBV	3410363	6840110	5.00%	1.001%
22	9.210	1287	1296	1304	rBV	364987	786730	0.58%	0.115%
23	9.463	1337	1339	1344	rVB4	3738	3601	0.00%	0.001%
24	9.522	1344	1349	1350	rBV3	5390	7609	0.01%	0.001%
25	9.698	1373	1379	1387	rVB3	19467	38301	0.03%	0.006%
26	9.851	1399	1405	1409	rBV3	22421	45052	0.03%	0.007%
27	9.910	1409	1415	1419	rVV2	80469	154808	0.11%	0.023%
28	9.981	1422	1427	1433	rVV	214502	400829	0.29%	0.059%
29	10.051	1433	1439	1452	rVB	97003	258174	0.19%	0.038%
30	10.163	1452	1458	1460	rBV6	12878	22791	0.02%	0.003%
31	10.269	1469	1476	1481	rBV	1130231	2091106	1.53%	0.306%
32	10.334	1481	1487	1495	rVB	2083800	3855621	2.82%	0.564%
33	10.457	1500	1508	1515	rVV2	401178	794161	0.58%	0.116%
34	10.528	1515	1520	1526	rVV	134388	252202	0.18%	0.037%
35	10.616	1529	1535	1542	rVV	312265	563584	0.41%	0.082%
36	10.710	1545	1551	1560	rVV2	219501	478556	0.35%	0.070%

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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\SFAMDLM110923SMA.M
 Title : SFAM01.0

37	10.816	1560	1569	1577	rVV	51553254	86493789	63.27%	12.660%
38	10.881	1577	1580	1588	rVB3	274484	521562	0.38%	0.076%
39	10.998	1594	1600	1606	rBV	71653	145268	0.11%	0.021%
40	11.063	1606	1611	1614	rBV2	149138	256106	0.19%	0.037%
41	11.128	1619	1622	1627	rVB	349529	551285	0.40%	0.081%
42	11.181	1627	1631	1637	rVB	162108	278028	0.20%	0.041%
43	11.239	1637	1641	1645	rVB	47564	66935	0.05%	0.010%
44	11.298	1645	1651	1655	rBV2	68596	118623	0.09%	0.017%
45	11.387	1655	1666	1676	rVB4	248567	737512	0.54%	0.108%
46	11.475	1676	1681	1692	rVB	116870	217463	0.16%	0.032%
47	11.610	1692	1704	1711	rBV	10580591	18528764	13.55%	2.712%
48	11.687	1711	1717	1727	rVB2	32889381	56812689	41.56%	8.316%
49	11.787	1727	1734	1735	rBV2	61026967	61437801	44.94%	8.993%
50	11.998	1767	1770	1779	rVB7	38722	86903	0.06%	0.013%
51	12.122	1780	1791	1793	rBV2	59775606	92958946	67.99%	13.606%
52	12.257	1812	1814	1816	rBV2	16684	18043	0.01%	0.003%
53	12.298	1817	1821	1824	rBV2	58767	93920	0.07%	0.014%
54	12.422	1836	1842	1851	rVB	1613975	2588456	1.89%	0.379%
55	12.581	1861	1869	1875	rVB	318587	596000	0.44%	0.087%
56	12.651	1875	1881	1888	rBV	242791	442712	0.32%	0.065%
57	12.763	1891	1900	1905	rVB	1052872	1668297	1.22%	0.244%
58	12.828	1905	1911	1914	rBV	3493146	5683542	4.16%	0.832%
59	12.904	1920	1924	1929	rVB	1599383	2304043	1.69%	0.337%
60	12.951	1929	1932	1936	rVB	262341	363381	0.27%	0.053%
61	12.992	1936	1939	1944	rVB2	82719	131735	0.10%	0.019%
62	13.063	1944	1951	1959	rBV	878376	1493626	1.09%	0.219%
63	13.134	1959	1963	1965	rBV2	59132	82083	0.06%	0.012%
64	13.216	1971	1977	1983	rBV	2947161	4577622	3.35%	0.670%
65	13.304	1987	1992	1997	rBV2	198659	296137	0.22%	0.043%
66	13.381	2001	2005	2007	rBV2	82201	113146	0.08%	0.017%
67	13.457	2012	2018	2024	rVV	3906185	6488128	4.75%	0.950%
68	13.539	2024	2032	2048	rVB	19542559	31911138	23.34%	4.671%
69	13.686	2052	2057	2058	rBV	145353	184495	0.13%	0.027%
70	13.716	2058	2062	2065	rVV2	513371	832169	0.61%	0.122%
71	13.757	2065	2069	2074	rVV2	687470	1296939	0.95%	0.190%
72	13.833	2075	2082	2090	rVV	10410777	17286120	12.64%	2.530%
73	13.916	2092	2096	2101	rVV	238088	338626	0.25%	0.050%
74	13.975	2101	2106	2109	rVV	475049	816405	0.60%	0.119%
75	14.004	2109	2111	2116	rVV	591943	856802	0.63%	0.125%
76	14.063	2116	2121	2127	rVB	1271876	1935279	1.42%	0.283%

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 Title : SFAM01.0

77	14.169	2134	2139	2144	rBV2	346681	569639	0.42%	0.083%
78	14.228	2145	2149	2151	rBV2	96083	153565	0.11%	0.022%
79	14.304	2157	2162	2166	rBV	346757	559201	0.41%	0.082%
80	14.386	2172	2176	2179	rBV	963408	1341155	0.98%	0.196%
81	14.428	2179	2183	2187	rVB	1588706	2283498	1.67%	0.334%
82	14.469	2187	2190	2194	rVB2	204863	234059	0.17%	0.034%
83	14.551	2194	2204	2209	rBV	3212040	5728378	4.19%	0.838%
84	14.598	2209	2212	2218	rVB4	475461	767427	0.56%	0.112%
85	14.657	2218	2222	2226	rBV2	292317	481356	0.35%	0.070%
86	14.698	2226	2229	2234	rVB	522485	718525	0.53%	0.105%
87	14.786	2235	2244	2249	rBV3	1164585	3104033	2.27%	0.454%
88	14.892	2256	2262	2266	rBV	772385	1307009	0.96%	0.191%
89	14.933	2266	2269	2278	rVB5	312434	527706	0.39%	0.077%
90	15.039	2282	2287	2290	rVV2	407631	686831	0.50%	0.101%
91	15.104	2291	2298	2311	rVB2	69179745	136714890	100.00%	20.011%
92	15.333	2327	2337	2343	rBV	462855	917484	0.67%	0.134%
93	15.445	2352	2356	2359	rBV2	71455	113621	0.08%	0.017%
94	15.528	2361	2370	2379	rVB	30572384	54908048	40.16%	8.037%
95	15.792	2410	2415	2419	rBV3	172095	317666	0.23%	0.046%
96	16.233	2483	2490	2499	rVB2	522299	1109920	0.81%	0.162%
97	16.339	2503	2508	2513	rVV4	106668	246853	0.18%	0.036%
98	16.404	2513	2519	2525	rVV	360422	813906	0.60%	0.119%
99	16.463	2526	2529	2535	rVV4	139122	252769	0.18%	0.037%
100	16.563	2536	2546	2566	rVB	3541334	8163113	5.97%	1.195%

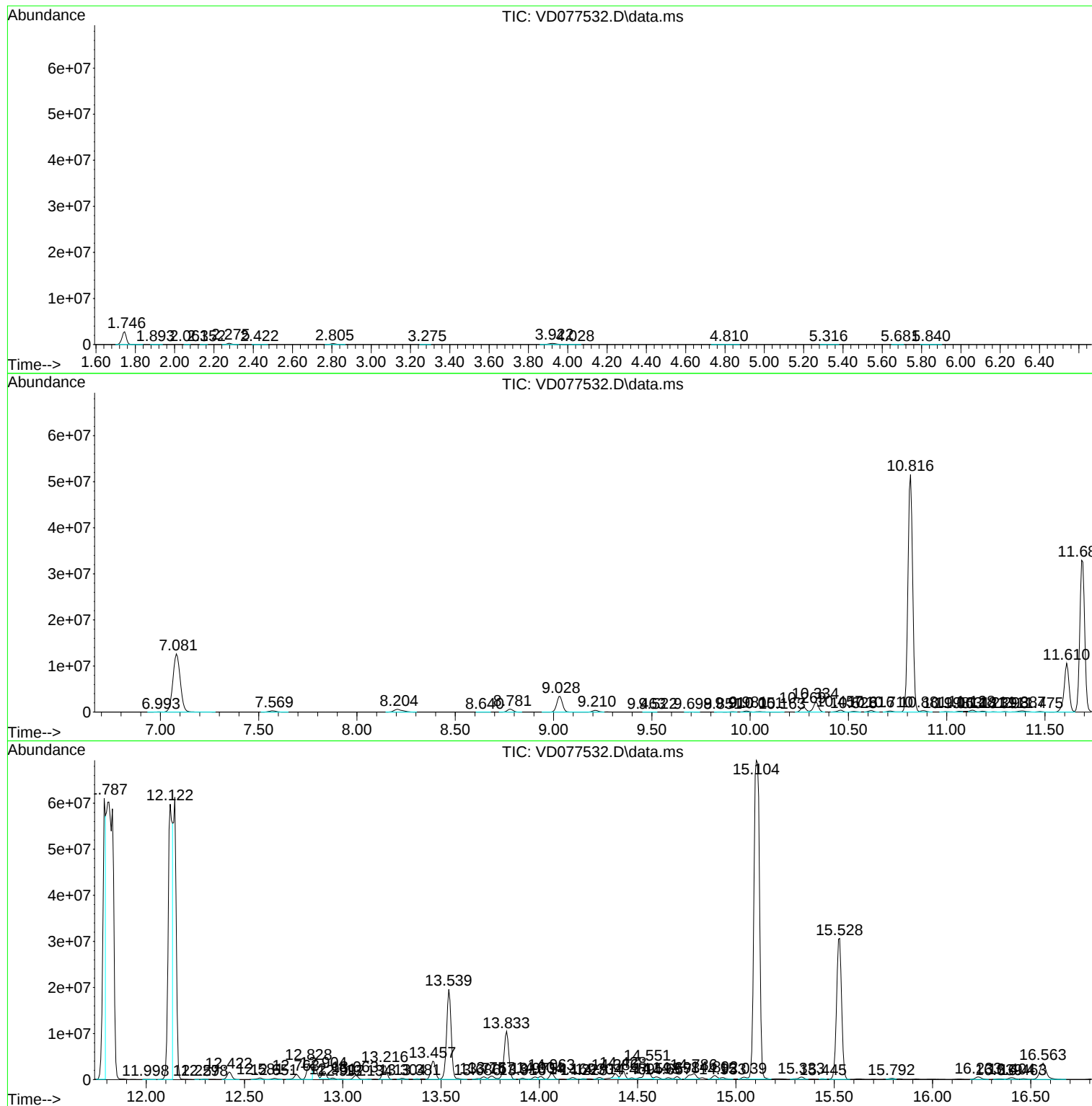
Sum of corrected areas: 683197091

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ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM110923SMA.M
Quant Title : SFAM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
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Sample : 05296-01
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Instrument :
MSVOA_D
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BH3R1

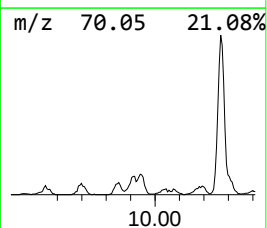
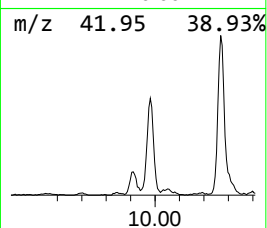
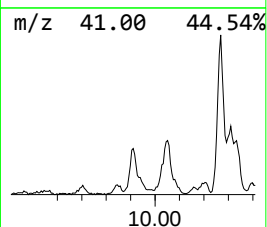
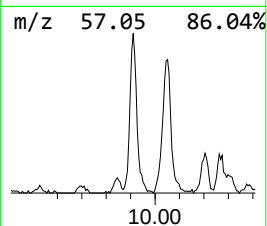
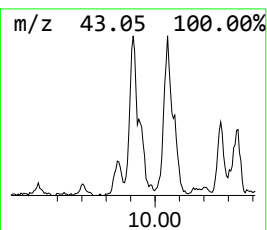
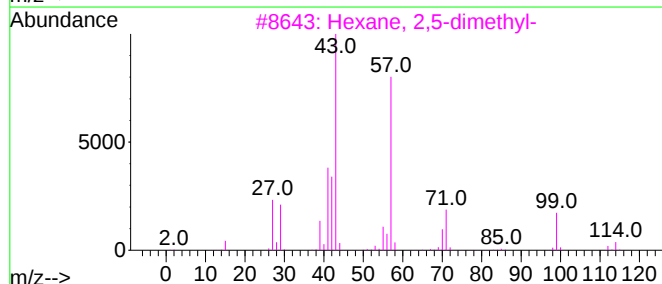
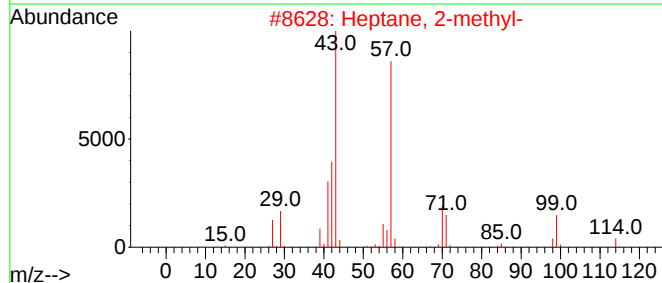
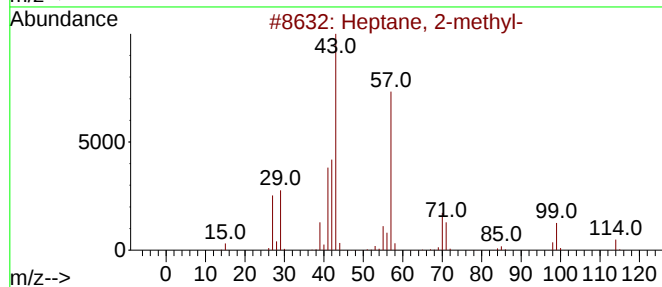
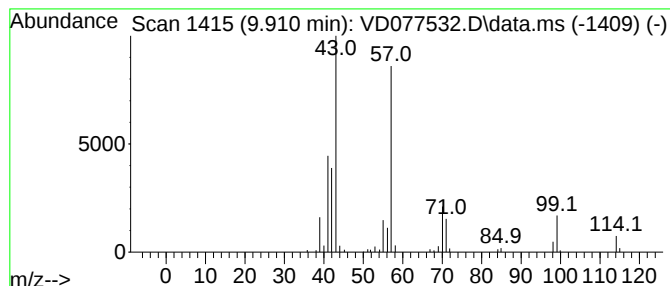
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM110923SMA.M
Quant Title : SFAM01.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.910	2.97 ug/L	154808	1,4-Difluorobenzene	8.781

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptane, 2-methyl-	114	C8H18	000592-27-8	94
2		Heptane, 2-methyl-	114	C8H18	000592-27-8	91
3		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	72
4		Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	58
5		Heptane, 2-methyl-	114	C8H18	000592-27-8	58



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Misc : 5.95G/10ml/MSVOA_D/SOIL
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Instrument :
MSVOA_D
ClientSampleId :
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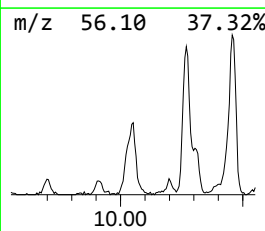
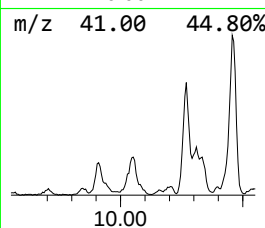
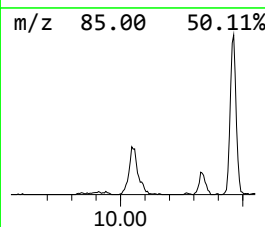
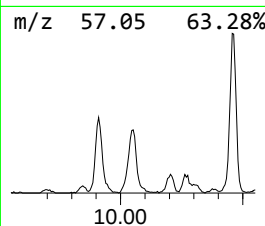
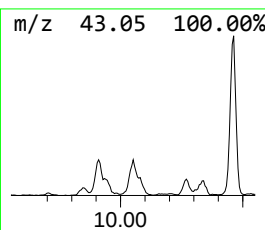
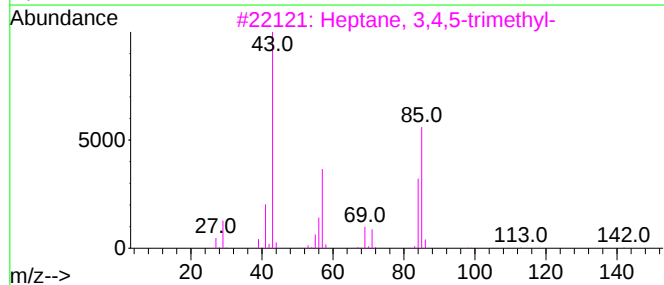
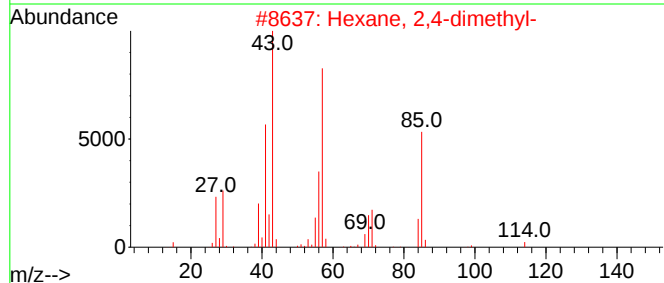
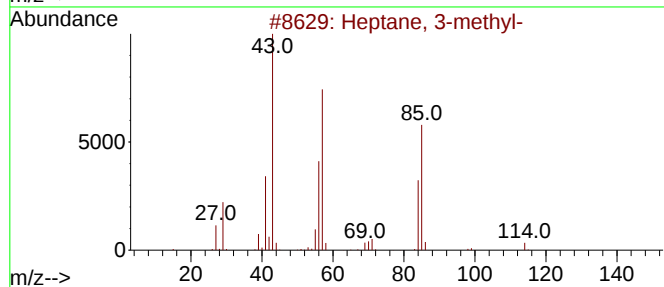
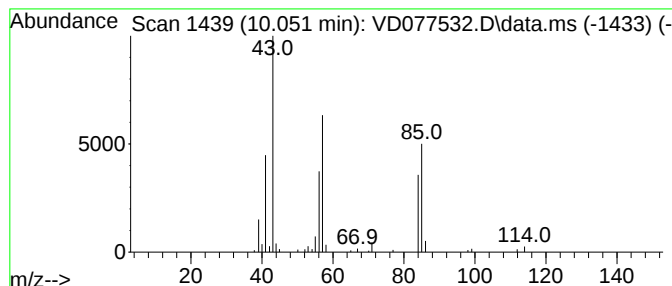
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM110923SMA.M
Quant Title : SFAM01.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.051	4.95 ug/L	258174	1,4-Difluorobenzene	8.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	80
3			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	78
4			Hexane, 2,3,3-trimethyl-	128	C9H20	016747-28-7	78
5			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72



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Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

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ClientSampleId :
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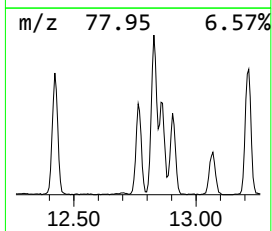
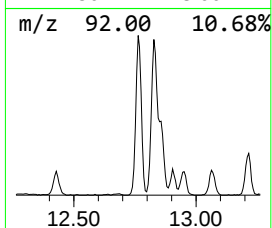
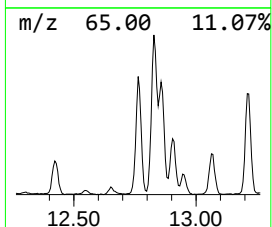
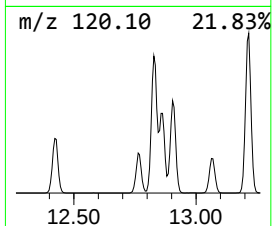
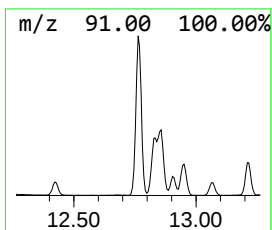
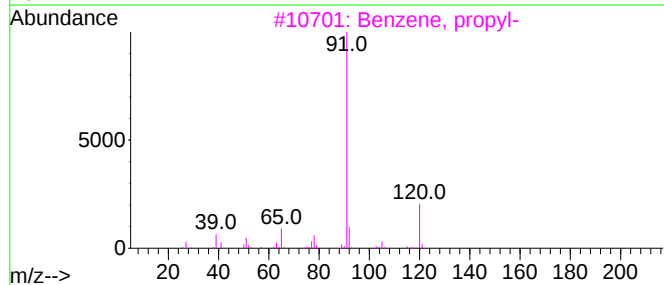
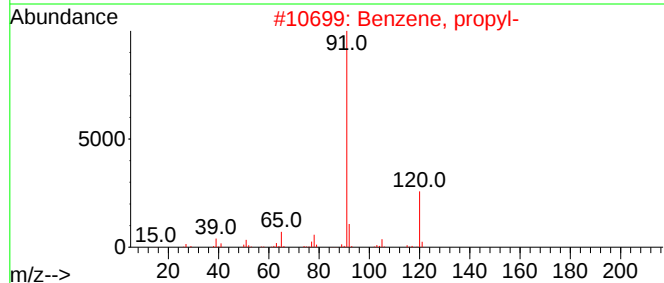
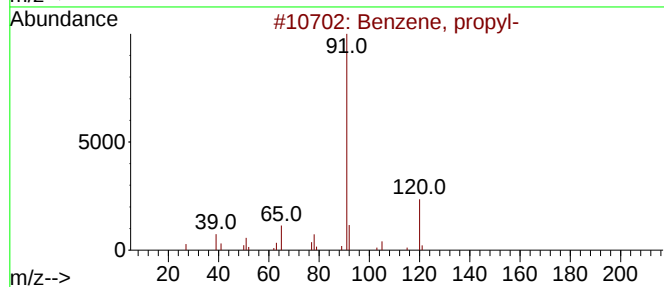
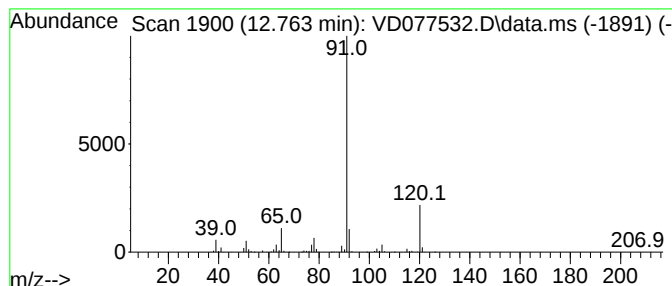
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM110923SMA.M
Quant Title : SFAM01.0

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

Peak Number 5 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.763	1.31 ug/L	1668300	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
5			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

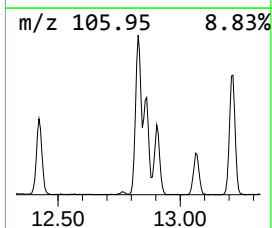
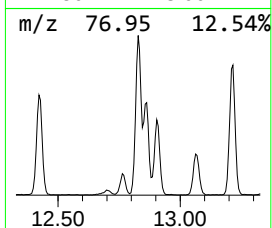
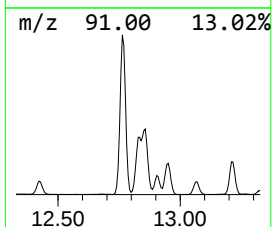
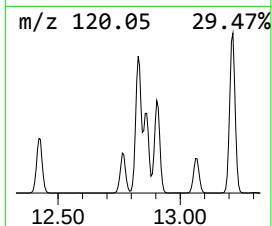
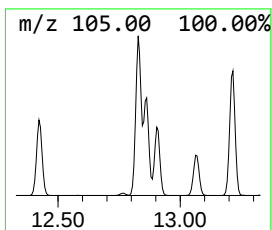
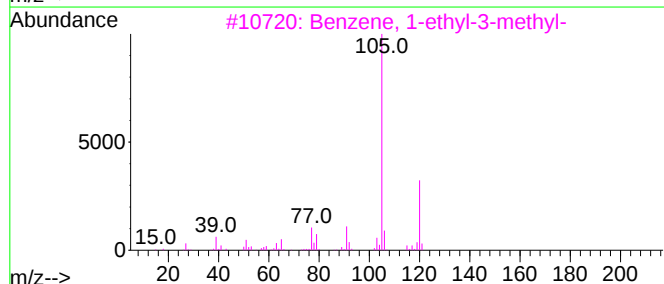
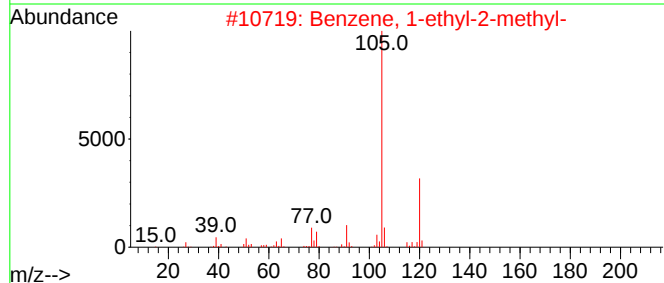
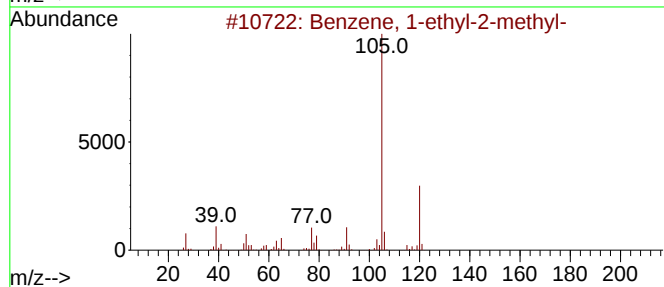
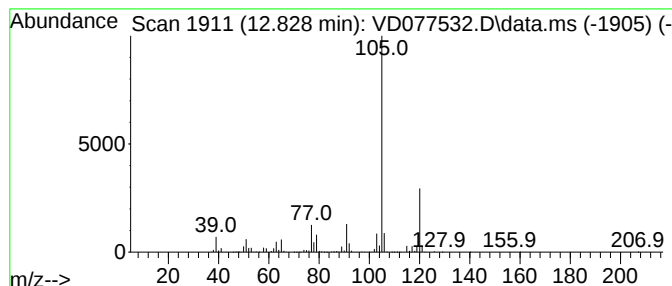
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Quant Title : SFAM01.0

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

Peak Number 6 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	4.45 ug/L	5683540	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

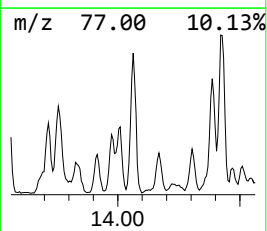
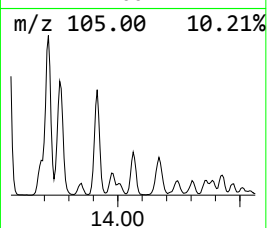
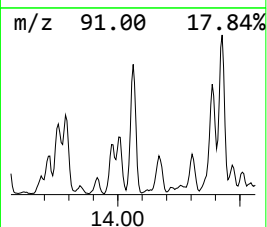
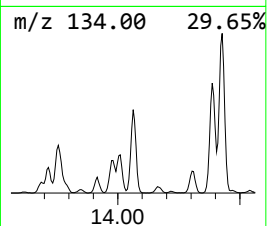
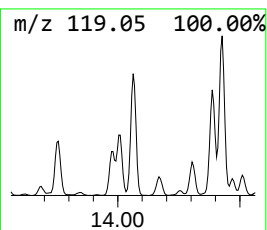
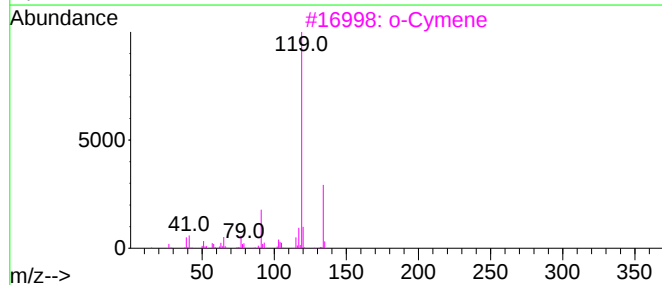
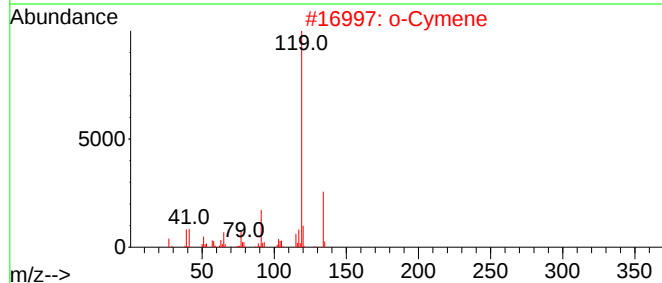
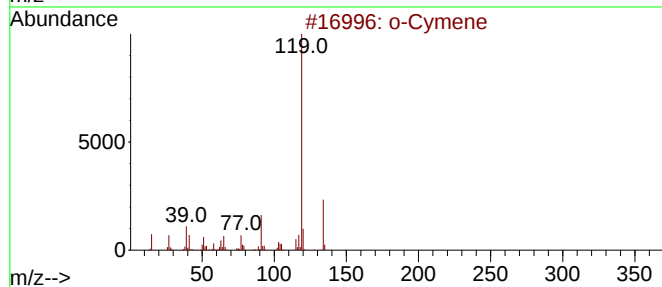
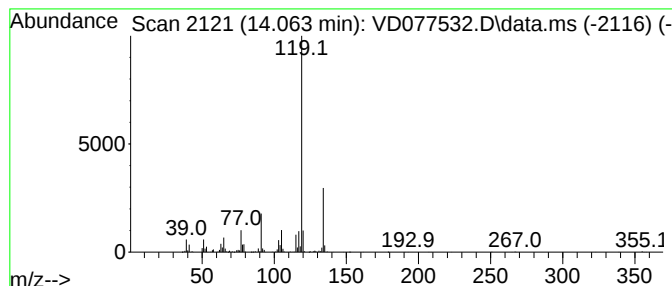
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Quant Title : SFAM01.0

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

Peak Number 7 o-Cymene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.063	1.52 ug/L	1935280	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			o-Cymene	134	C10H14	000527-84-4	95
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

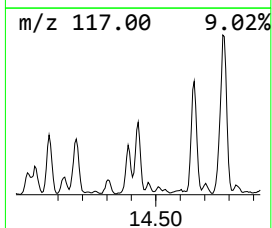
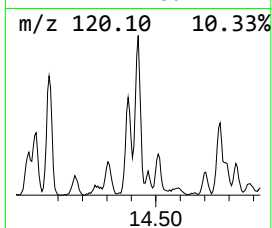
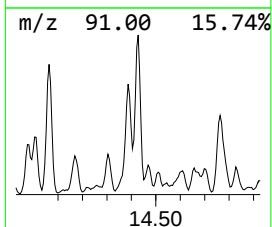
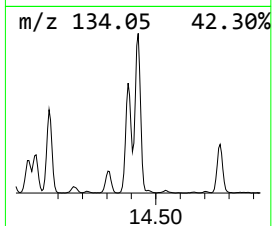
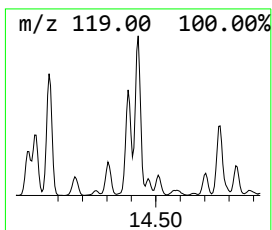
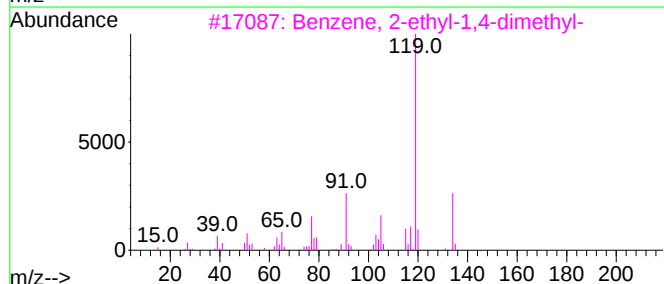
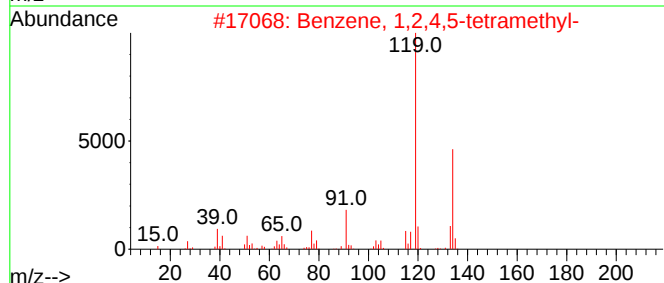
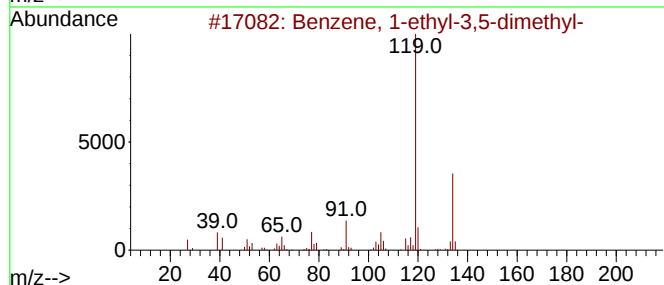
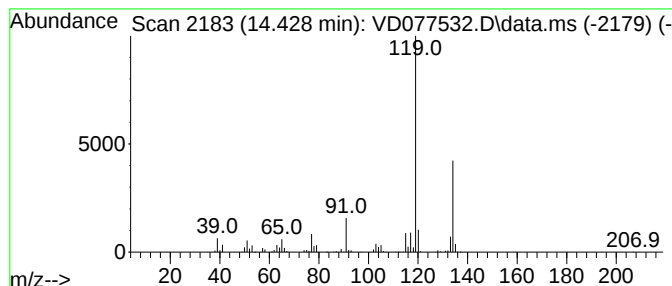
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Quant Title : SFAM01.0

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

Peak Number 8 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.428	1.79 ug/L	2283500	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			o-Cymene	134	C10H14	000527-84-4	94
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

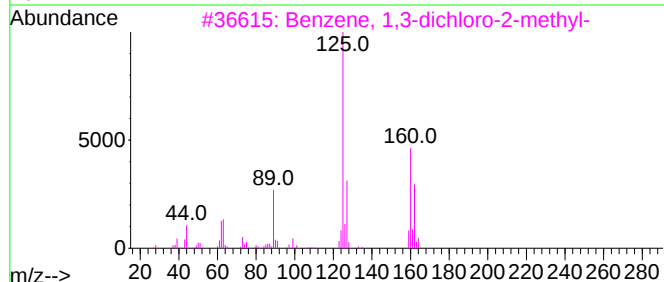
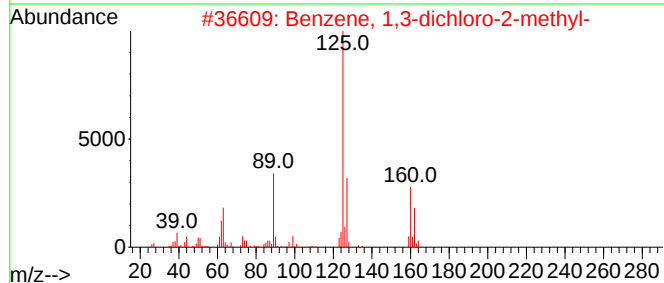
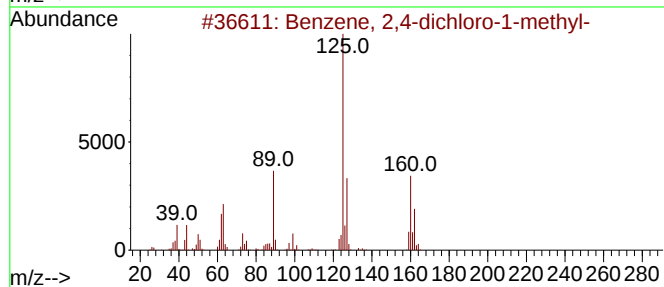
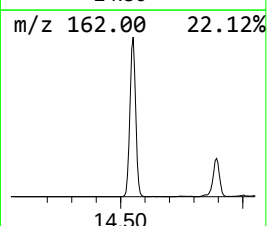
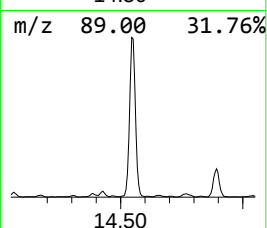
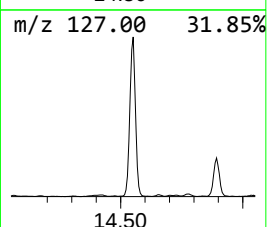
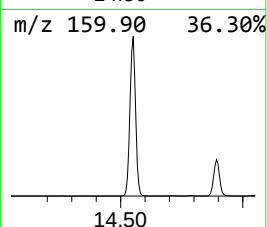
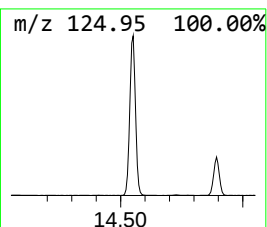
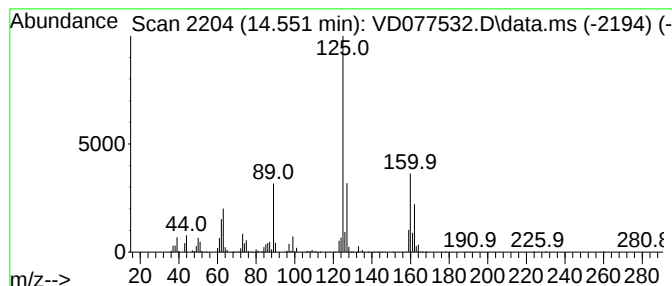
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Quant Title : SFAM01.0

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

Peak Number 9 Benzene, 2,4-dichloro-1-met... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.551	4.49 ug/L	5728380	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2,4-dichloro-1-methyl-	160	C7H6Cl2	000095-73-8	98
2			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
3			Benzene, 1,3-dichloro-2-methyl-	160	C7H6Cl2	000118-69-4	97
4			Benzene, 1,2-dichloro-4-methyl-	160	C7H6Cl2	000095-75-0	97
5			Benzene, 1,2-dichloro-3-methyl-	160	C7H6Cl2	032768-54-0	96



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

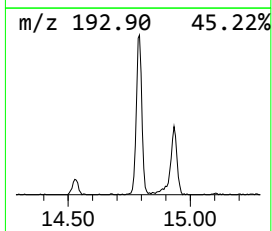
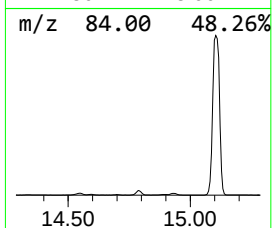
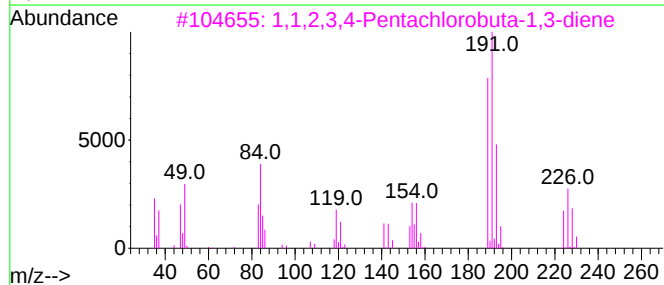
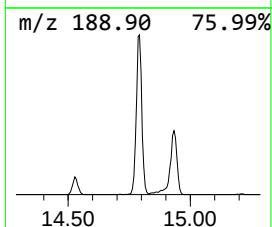
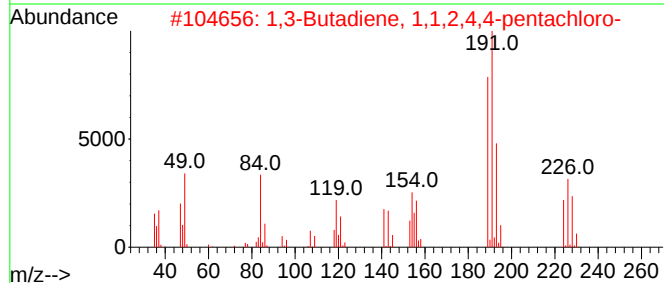
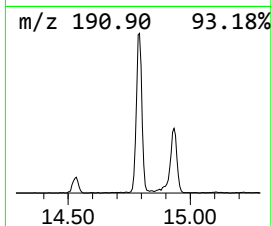
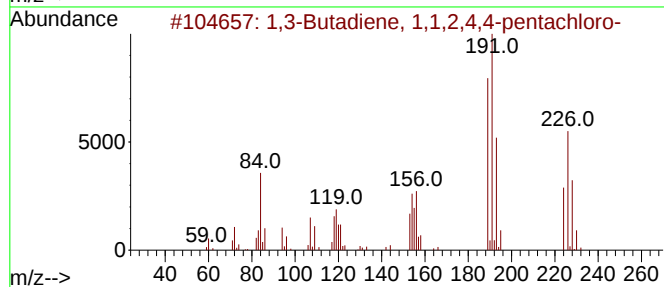
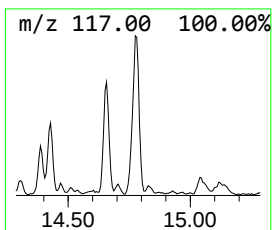
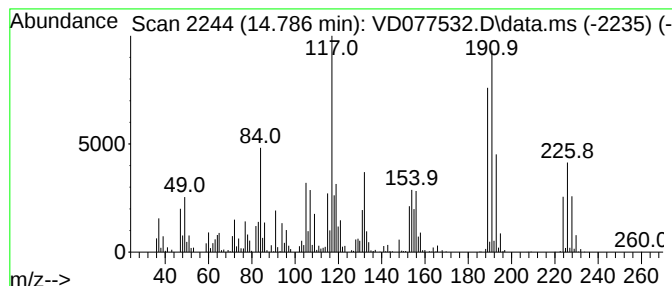
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Quant Title : SFAM01.0

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

Peak Number 10 1,3-Butadiene, 1,1,2,4,4-pe... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.786	2.43 ug/L	3104030	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Butadiene, 1,1,2,4,4-pentach...	224	C4HCl5	021400-41-9	99		
2	1,3-Butadiene, 1,1,2,4,4-pentach...	224	C4HCl5	021400-41-9	60		
3	1,1,2,3,4-Pentachlorobuta-1,3-diene	224	C4HCl5	1010487-10-4	38		
4	2-Thiophenecarbonyl chloride, 5-...	224	C5H2BrClO5	031555-60-9	27		
5	2-Methoxy-5-ethyl-5-phenyl-4-oxa...	221	C12H15NO3	056440-40-5	16		



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

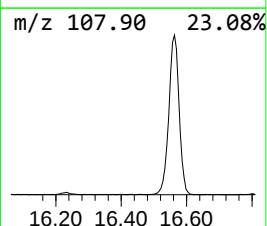
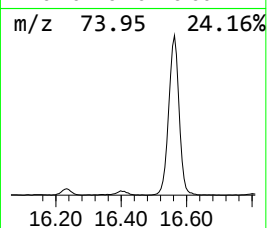
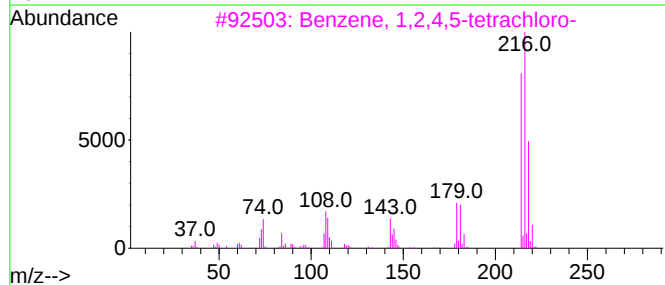
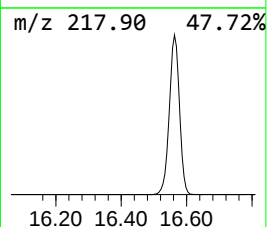
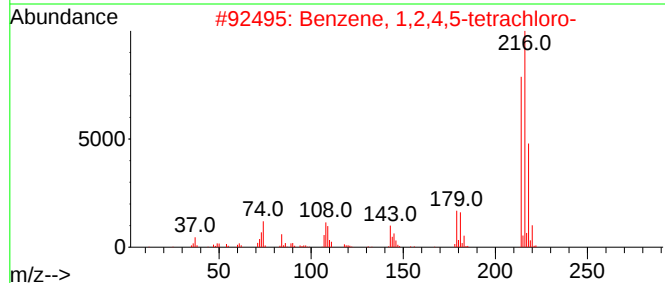
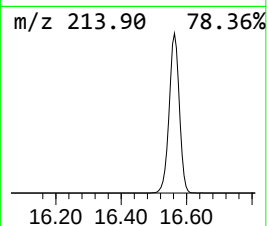
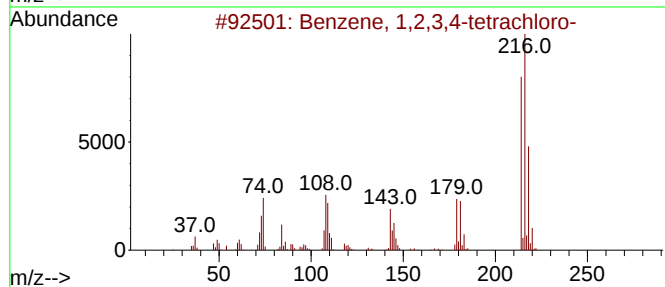
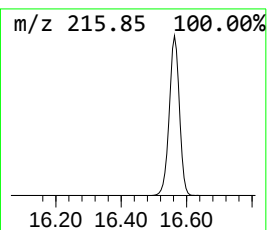
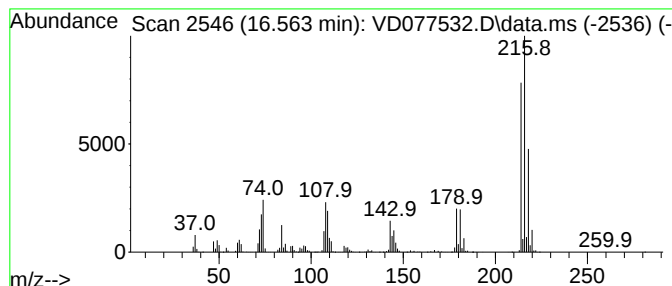
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Quant Title : SFAM01.0

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

Peak Number 11 Benzene, 1,2,3,4-tetrachloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.563	6.40 ug/L	8163110	1,4-Dichlorobenzene-d4	13.522

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	99
2			Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	99
3			Benzene, 1,2,4,5-tetrachloro-	214	C6H2Cl4	000095-94-3	99
4			Benzene, 1,2,3,5-tetrachloro-	214	C6H2Cl4	000634-90-2	99
5			Benzene, 1,2,3,4-tetrachloro-	214	C6H2Cl4	000634-66-2	99



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD110923\
Data File : VD077532.D
Acq On : 09 Nov 2023 13:30
Operator : JC/SY
Sample : 05296-01
Misc : 5.95G/10ml/MSVOA_D/SOIL
ALS Vial : 1 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
BH3R1

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\SFAMDLM110923SMA.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
(DEL) Alkane: S...	9.910	3.0	ug/L	154808	1	8.781	1302920	25.0
(DEL) Alkane: S...	10.051	5.0	ug/L	258174	1	8.781	1302920	25.0
Benzene, propyl-	12.763	1.3	ug/L	1668300	3	13.522	31911100	25.0
Benzene, 1-ethy...	12.828	4.5	ug/L	5683540	3	13.522	31911100	25.0
o-Cymene	14.063	1.5	ug/L	1935280	3	13.522	31911100	25.0
Benzene, 1-ethy...	14.428	1.8	ug/L	2283500	3	13.522	31911100	25.0
Benzene, 2,4-di...	14.551	4.5	ug/L	5728380	3	13.522	31911100	25.0
1,3-Butadiene, ...	14.786	2.4	ug/L	3104030	3	13.522	31911100	25.0
Benzene, 1,2,3,...	16.563	6.4	ug/L	8163110	3	13.522	31911100	25.0