LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_X\Method\82X053123W.M
Title : SW846 8260

Signal : TIC: VX036010.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.142	6	9	22	rVB	653648	837720	24.72%	1.974%
2	1.264	26	29	40	rBV	205620	198640	5.86%	0.468%
3	1.367	42	46	53	rBV	434194	423206	12.49%	0.997%
4	1.751	103	109	123	rBV	2004242	2676874	79.00%	6.309%
5	1.953	137	142	156	rVB	343767	480490	14.18%	1.132%
6	2.068	156	161	166	rBV	24879	36144	1.07%	0.085%
7	2.215	180	185	197	rVB	105658	160610	4.74%	0.379%
8	2.343	197	206	212	rBV	170325	337396	9.96%	0.795%
9	2.782	259	278	281	rBV	283914	655910	19.36%	1.546%
10	2.824	281	285	289	rVV	479729	982821	29.00%	2.316%
11	2.867	289	292	308	rVB2	249818	525700	15.51%	1.239%
12	3.099	320	330	345	rVB	443203	992538	29.29%	2.339%
13	3.446	378	387	398	rVB2	27491	61526	1.82%	0.145%
14	3.727	426	433	443	rVB	37710	90617	2.67%	0.214%
15	3.836	443	451	457	rBV2	36679	92352	2.73%	0.218%
16	3.904	457	462	470	rVB2	15318	37222	1.10%	0.088%
17	4.105	488	495	502	rVB2	19816	48809	1.44%	0.115%
18	4.208	504	512	518	rVV2	27446	78509	2.32%	0.185%
19	4.300	518	527	540	rVV	569773	1641988	48.46%	3.870%
20	4.428	540	548	564	rVB6	17328	61737	1.82%	0.146%
21	5.019	636	645	654	rBV3	11848	39039	1.15%	0.092%
22	5.385	690	705	711	rBV	138949	400231	11.81%	0.943%
23	5.470	711	719	726	rVV	260370	840460	24.80%	1.981%
24	5.550	726	732	739	rVV2	278784	836005	24.67%	1.970%
25	5.623	739	744	765	rVB2	115753	364655	10.76%	0.859%
26	5.830	765	778	789	rBV5	53125	172576	5.09%	0.407%
27	5.958	789	799	806	rVV	159925	436817	12.89%	1.029%
28	6.037	806	812	822	rVV	135756	361403	10.67%	0.852%
29	6.159	822	832	841	rVV6	77112	320797	9.47%	0.756%
30	6.257	841	848	857	rVV4	122483	376173	11.10%	0.887%
31	6.354	857	864	875	rVB2	57564	164608	4.86%	0.388%
32	6.604	897	905	921	rVB6	9950	37032	1.09%	0.087%
33	6.763	921	931	939	rBV	392065	953565	28.14%	2.247%
34	6.842	939	944	956	rVB4	28517	83043	2.45%	0.196%
35	7.171	991	998	1006	rBV3	22454	52121	1.54%	0.123%
36	7.372	1017	1031	1042	rBV4	200677	569430	16.81%	1.342%

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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_X\Method\82X053123W.M
Title : SW846 8260

37	7.659	1072	1078	1086	rVB	43154	85693	2.53%	0.202%
38	7.854	1103	1110	1112	rBV2	31620	59265	1.75%	0.140%
39	7.885	1112	1115	1122	rVB2	35614	61201	1.81%	0.144%
40	7.982	1122	1131	1144	rVB	42410	97408	2.87%	0.230%
41	8.110	1144	1152	1154	rBV	111974	220148	6.50%	0.519%
42	8.141	1154	1157	1163	rVB	104380	168276	4.97%	0.397%
43	8.427	1199	1204	1211	rBV	19506	34034	1.00%	0.080%
44	8.506	1211	1217	1225	rVB3	75150	133911	3.95%	0.316%
45	8.604	1225	1233	1235	rBV4	22685	47684	1.41%	0.112%
46	8.647	1235	1240	1247	rVV	792073	1233224	36.39%	2.906%
47	8.720	1247	1252	1257	rVB	96877	146517	4.32%	0.345%
48	8.842	1264	1272	1280	rVB2	48413	98923	2.92%	0.233%
49	8.957	1286	1291	1299	rVB	237218	382430	11.29%	0.901%
50	9.049	1299	1306	1311	rVV	233063	374557	11.05%	0.883%
51	9.098	1311	1314	1319	rVV	28407	46060	1.36%	0.109%
52	9.159	1319	1324	1335	rVB	60321	99362	2.93%	0.234%
53	9.451	1364	1372	1376	rBV4	17533	41371	1.22%	0.098%
54	10.055	1460	1471	1476	rBV	805479	1104940	32.61%	2.604%
55	10.146	1481	1486	1490	rVV2	42433	70441	2.08%	0.166%
56	10.195	1490	1494	1499	rVB	78419	97767	2.89%	0.230%
57	10.299	1505	1511	1517	rVB	486165	676556	19.97%	1.595%
58	10.640	1563	1567	1575	rVB	49345	66723	1.97%	0.157%
59	10.963	1614	1620	1625	rBV3	31156	46656	1.38%	0.110%
60	11.079	1634	1639	1645	rBV	646530	816434	24.09%	1.924%
61	11.305	1672	1676	1684	rVB2	43832	62503	1.84%	0.147%
62	11.396	1684	1691	1696	rVV	126826	188087	5.55%	0.443%
63	11.451	1696	1700	1710	rVB	41275	61644	1.82%	0.145%
64	11.609	1721	1726	1735	rVB	803203	1009412	29.79%	2.379%
65	11.750	1744	1749	1755	rBV	1433167	1695411	50.03%	3.996%
66	11.969	1778	1785	1788	rBV	30131	38445	1.13%	0.091%
67	12.018	1788	1793	1799	rVB	850577	1096167	32.35%	2.583%
68	12.085	1799	1804	1810	rBV	237815	291593	8.61%	0.687%
69	12.176	1815	1819	1824	rVB	38456	46740	1.38%	0.110%
70	12.243	1824	1830	1836	rBV	2734801	3388455	100.00%	7.986%
71	12.311	1836	1841	1848	rVB	338029	409139	12.07%	0.964%
72	12.402	1851	1856	1861	rVB2	141425	179280	5.29%	0.423%
73	12.463	1861	1866	1871	rVB	241240	297242	8.77%	0.701%
74	12.530	1871	1877	1885	rBV	303217	457382	13.50%	1.078%
75	12.609	1885	1890	1894	rBV	723698	929505	27.43%	2.191%
76	12.646	1894	1896	1900	rVB	242479	245426	7.24%	0.578%

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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Title : SW846 8260

77	12.701	1900	1905	1912	rBV2	876166	1253037	36.98%	2.953%
78	12.847	1923	1929	1934	rVB2	103119	143198	4.23%	0.337%
79	12.920	1934	1941	1945	rBV	942060	1122193	33.12%	2.645%
80	12.963	1945	1948	1954	rVB	342552	401410	11.85%	0.946%
81	13.024	1954	1958	1966	rVB2	64829	100040	2.95%	0.236%
82	13.134	1966	1976	1978	rBV3	57048	100863	2.98%	0.238%
83	13.170	1978	1982	1992	rVB	963845	1187528	35.05%	2.799%
84	13.292	1992	2002	2007	rBV2	2001991	2639933	77.91%	6.222%
85	13.341	2007	2010	2014	rVB3	42888	51568	1.52%	0.122%
86	13.420	2019	2023	2029	rVB	170433	222165	6.56%	0.524%
87	13.505	2031	2037	2039	rBV2	70319	88596	2.61%	0.209%
88	13.542	2039	2043	2046	rVV	167464	229257	6.77%	0.540%
89	13.585	2046	2050	2054	rVV3	178011	290213	8.56%	0.684%
90	13.664	2055	2063	2070	rVB2	195599	427262	12.61%	1.007%
91	13.774	2074	2081	2090	rVB	151863	211226	6.23%	0.498%
92	13.926	2100	2106	2110	rBV2	72403	103489	3.05%	0.244%
93	13.969	2110	2113	2117	rVB	74215	83967	2.48%	0.198%
94	14.072	2125	2130	2136	rBV	120867	154655	4.56%	0.364%
95	14.213	2148	2153	2158	rBV	91810	140763	4.15%	0.332%
96	14.316	2167	2170	2175	rBV	46032	60867	1.80%	0.143%
97	14.371	2175	2179	2184	rVB	78253	93081	2.75%	0.219%
98	14.597	2209	2216	2220	rBV	30768	45976	1.36%	0.108%
99	14.780	2241	2246	2255	rBV	185162	244283	7.21%	0.576%

Sum of corrected areas: 42430346

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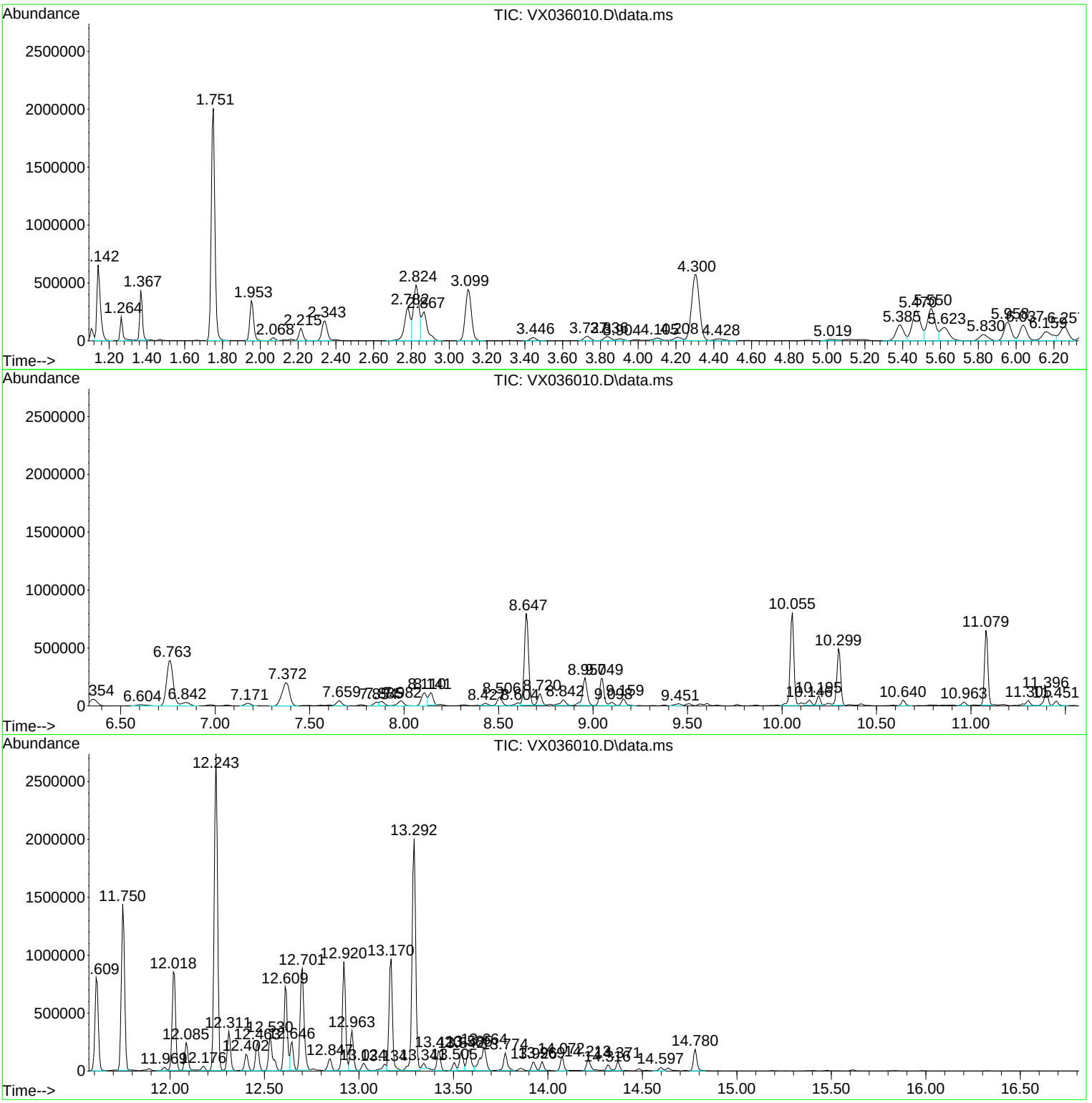
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
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ClientSampleId :
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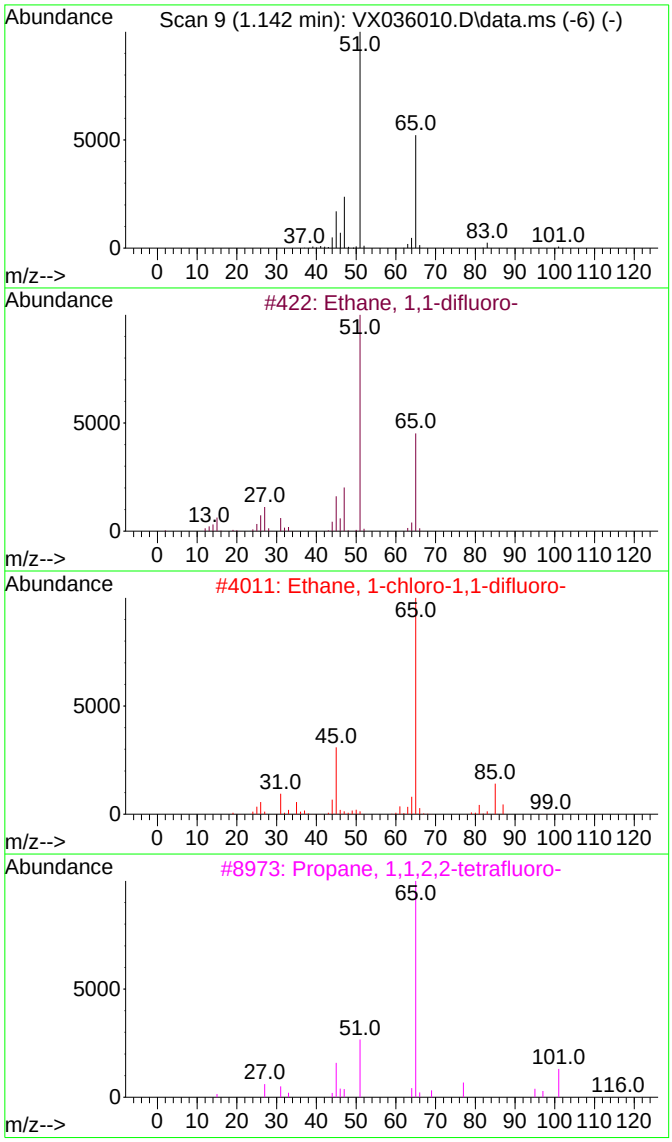
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

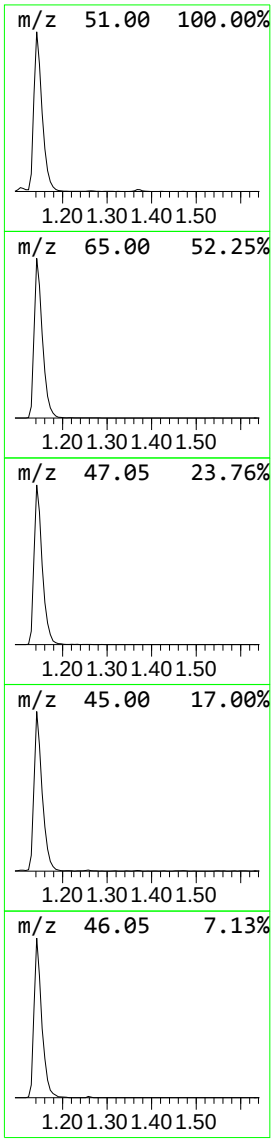
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Ethane, 1,1-difluoro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.142	50.10 ug/l	837720	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Ethane, 1,1-difluoro-	66	C2H4F2	000075-37-6	91
2	Ethane, 1-chloro-1,1-difluoro-	100	C2H3ClF2	000075-68-3	9
3	Propane, 1,1,2,2-tetrafluoro-	116	C3H4F4	040723-63-5	4
4	Difluorochloromethane	86	CHClF2	000075-45-6	4
5	Ethanamine, N-chloro-N,1,1-trifl...	133	C2H3ClF3N	016276-45-2	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

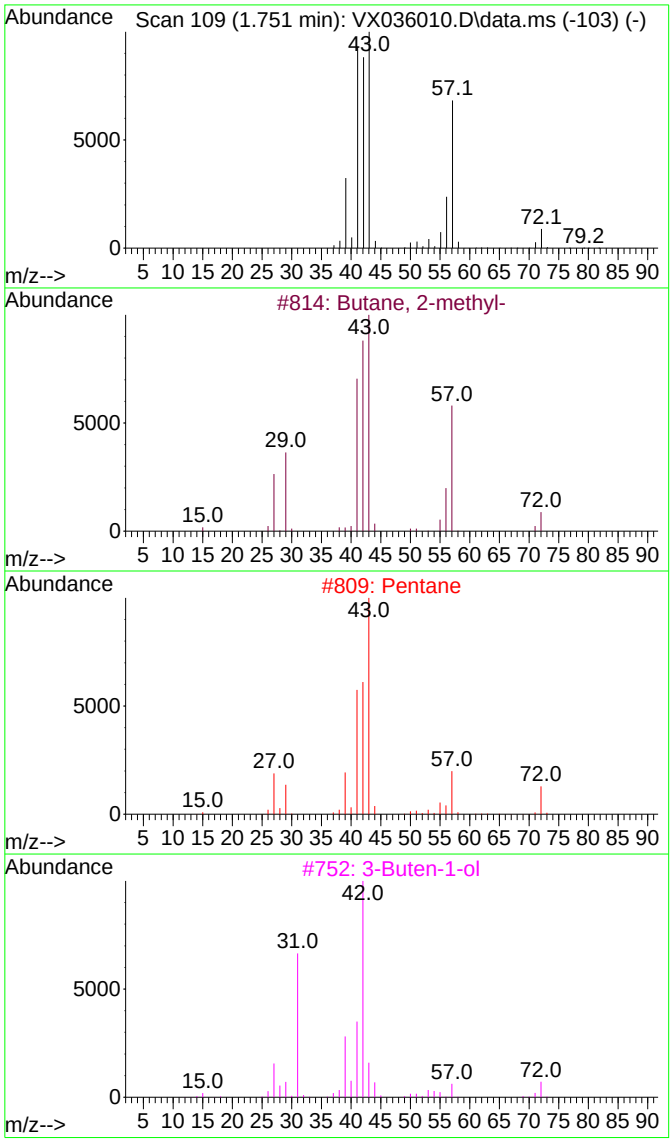
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

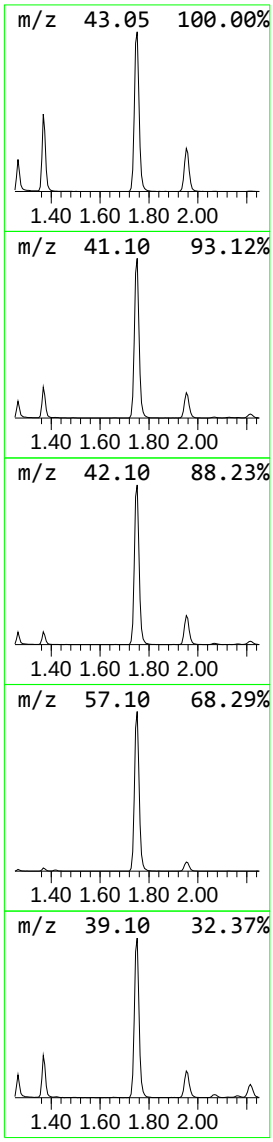
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.751	160.10 ug/l	2676870	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Pentane	72	C5H12	000109-66-0	36
3	3-Buten-1-ol	72	C4H8O	000627-27-0	28
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5	Methane, isocyanato-	57	C2H3NO	000624-83-9	9





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MSVOA_X
ClientSampleId :
MH-289

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

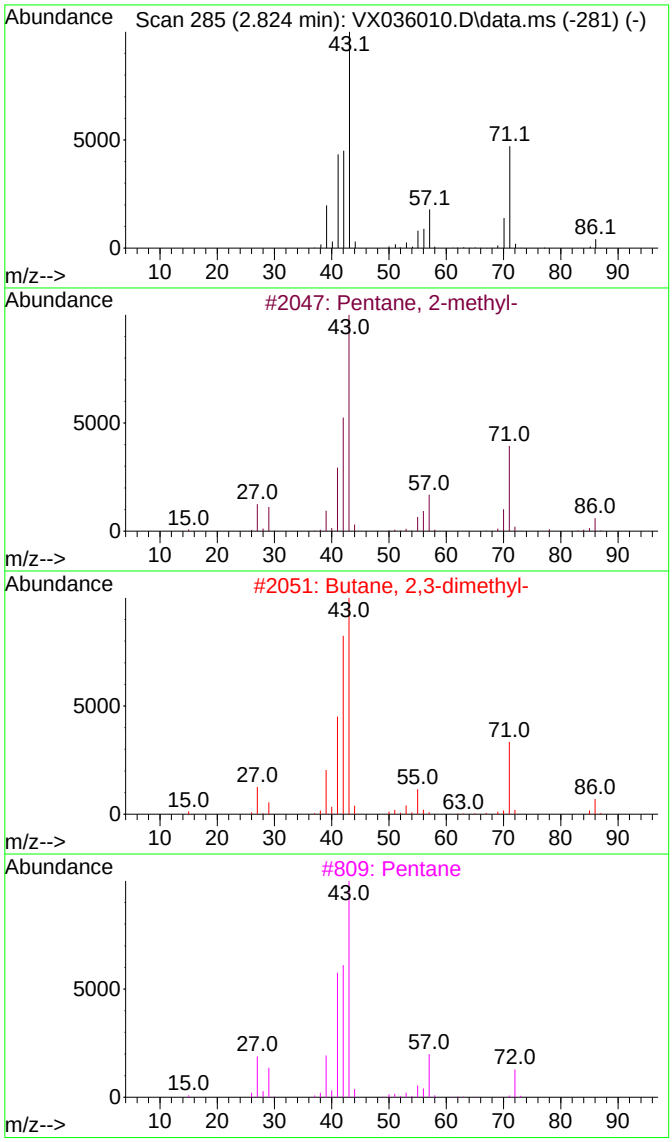
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

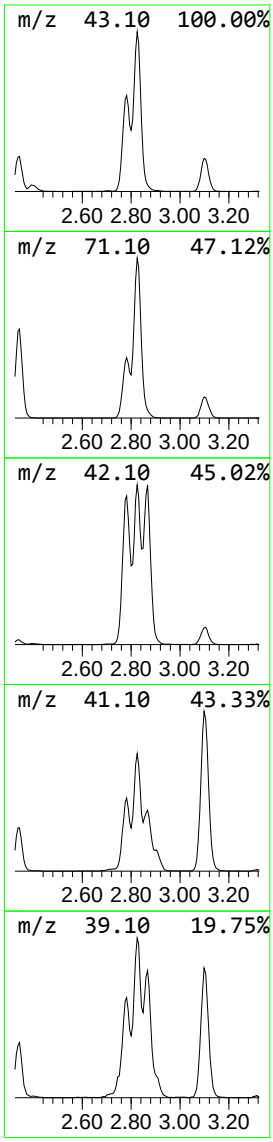
****

Peak Number 3 Pentane, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.824	58.78 ug/l	982821	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	47
3	Pentane	72	C5H12	000109-66-0	47
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	36
5	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	28





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Instrument :
MSVOA_X
ClientSampleId :
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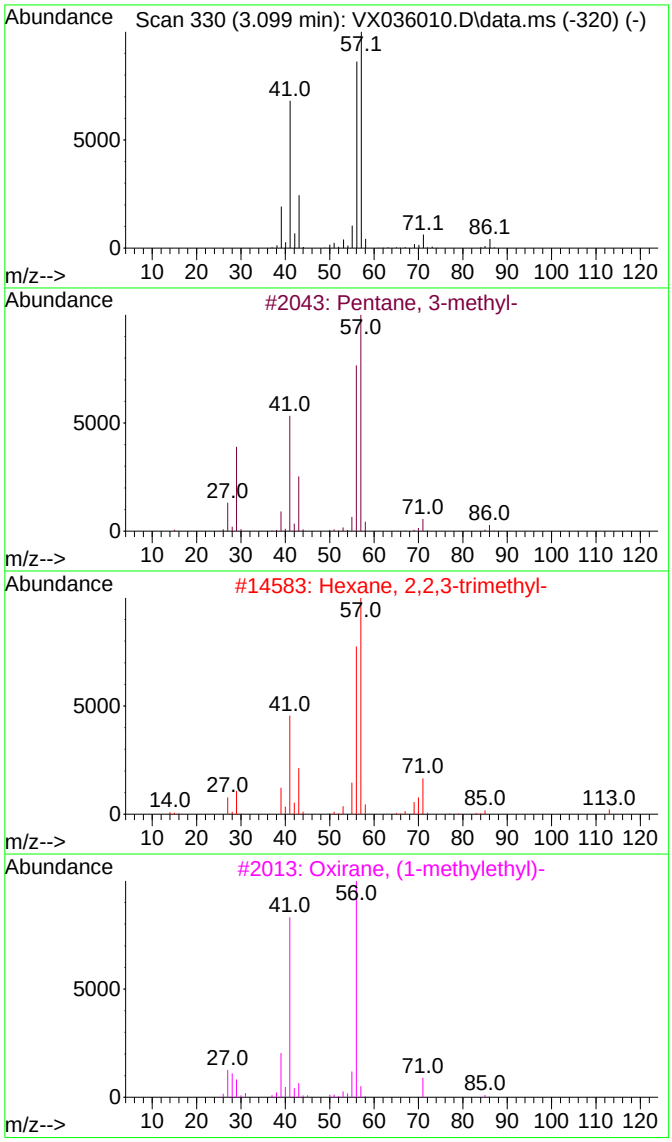
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

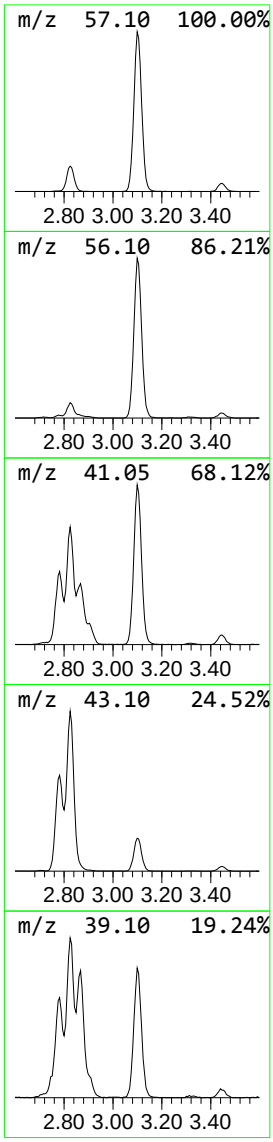
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.099	59.36 ug/l	992538	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43
5	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	42





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Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

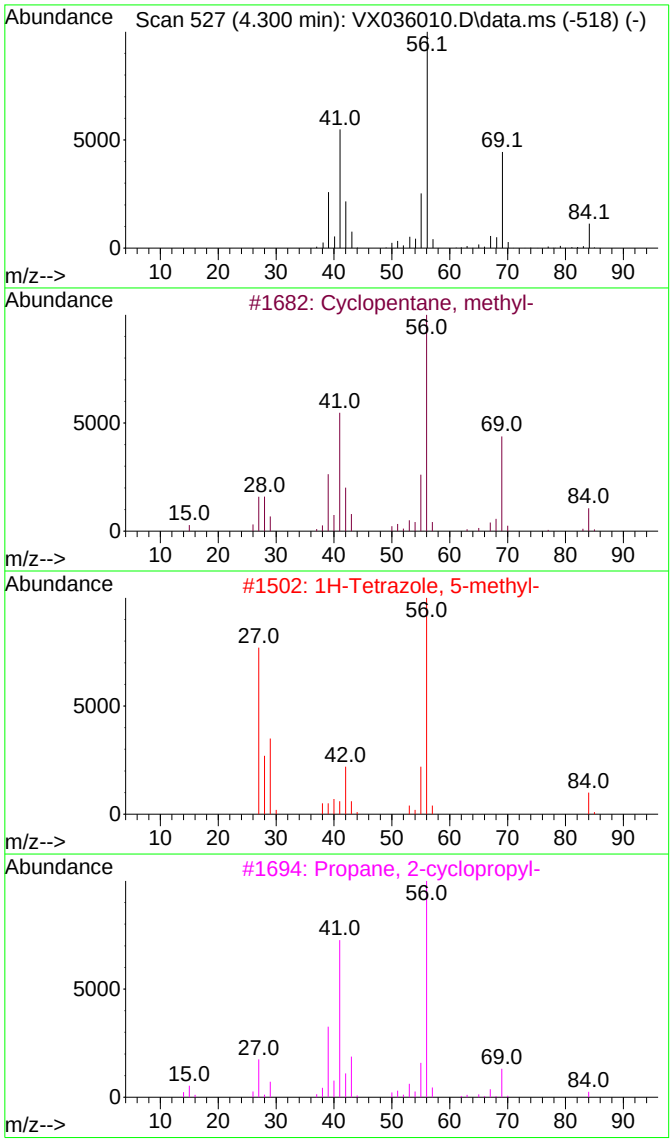
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

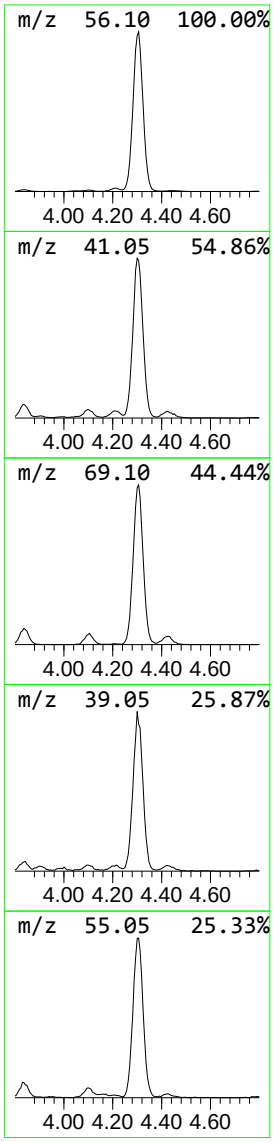
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.300	98.20 ug/l	1641990	Pentafluorobenzene	5.556

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
3	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	72
4	Cyclohexane	84	C6H12	000110-82-7	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

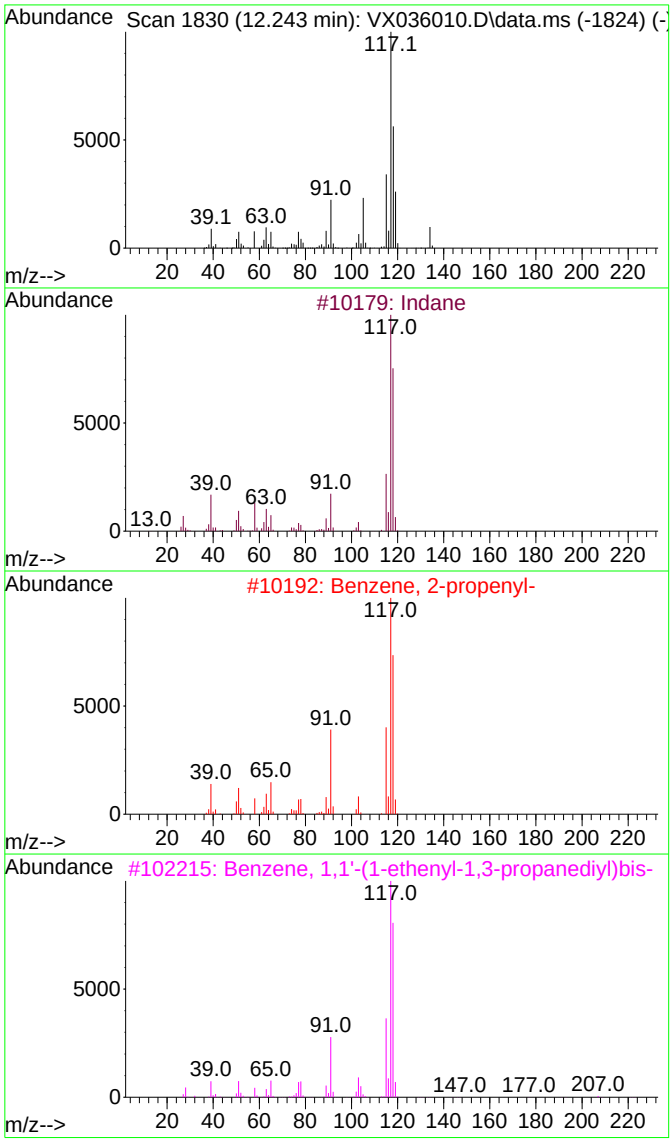
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

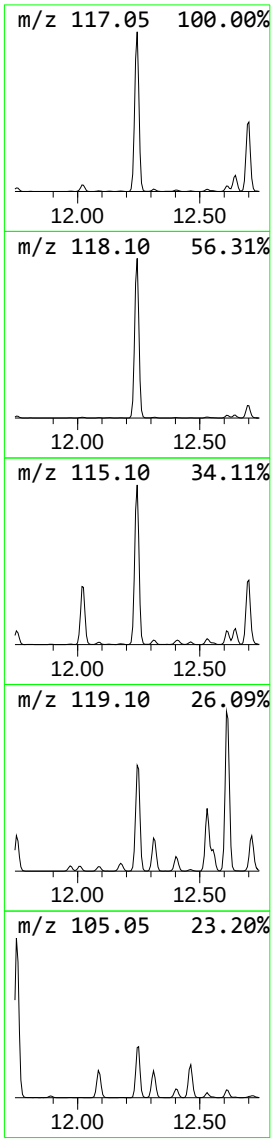
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.243	154.56 ug/l	3388460	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	70
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	70
3	Benzene, 1,1'-(1-ethenyl-1,3-pro...	222	C17H18	061141-97-7	64
4	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	64
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

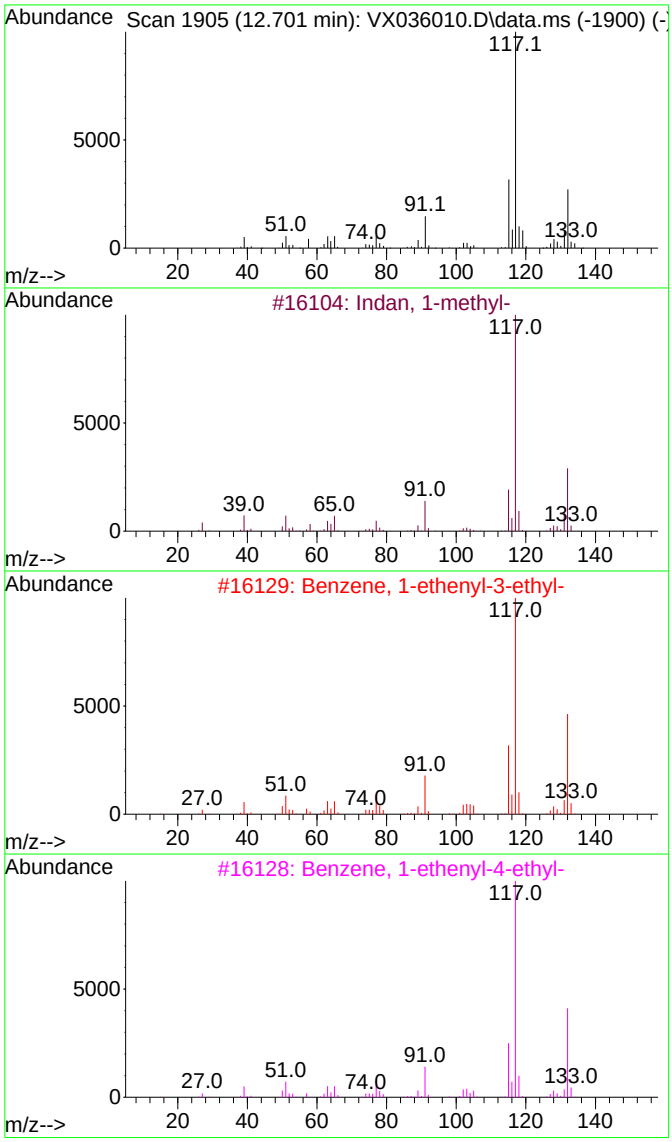
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

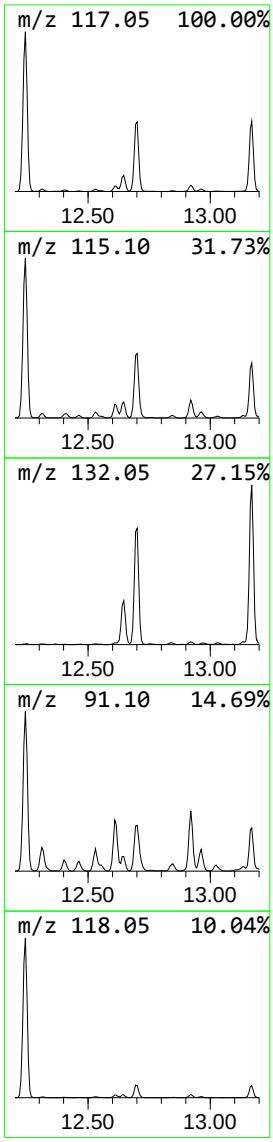
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.701	57.16 ug/l	1253040	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

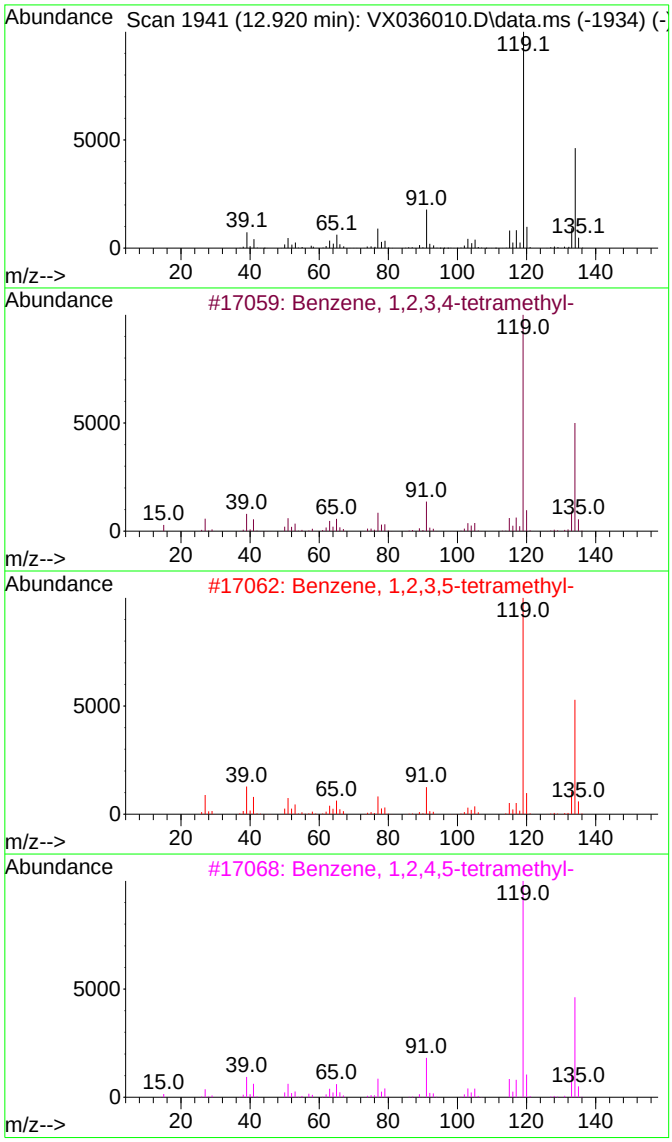
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

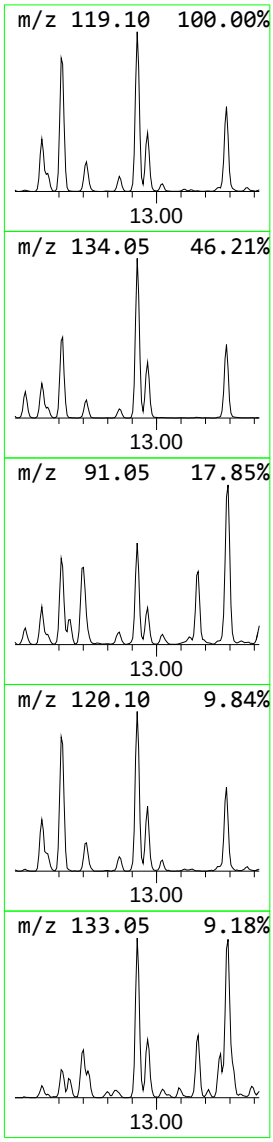
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	51.19 ug/l	1122190	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

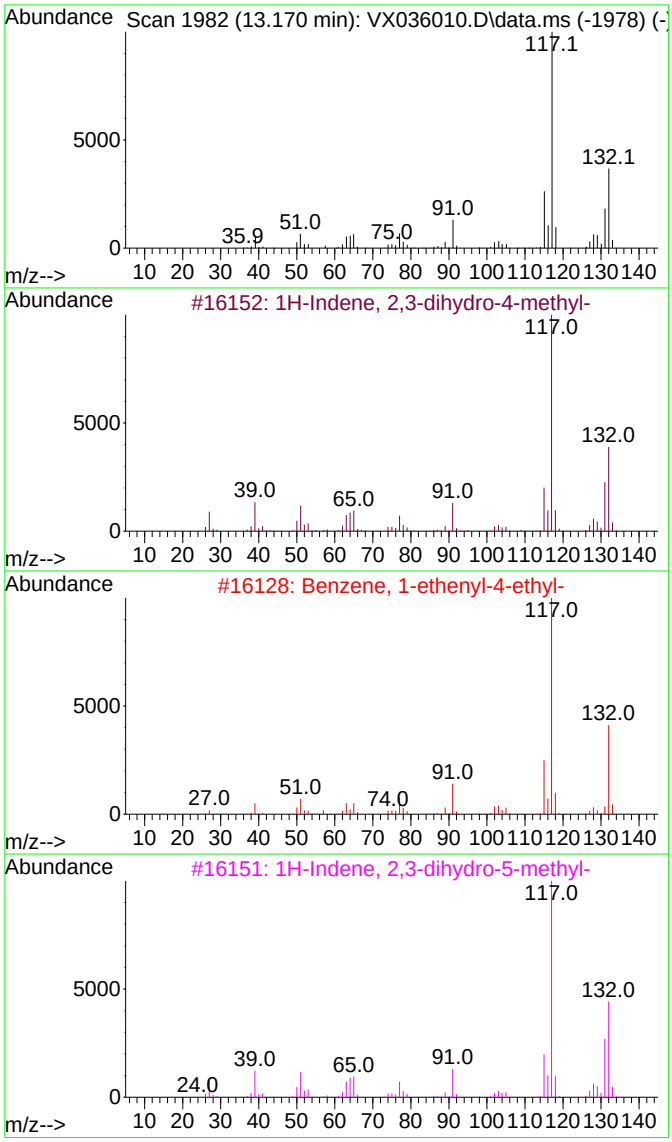
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

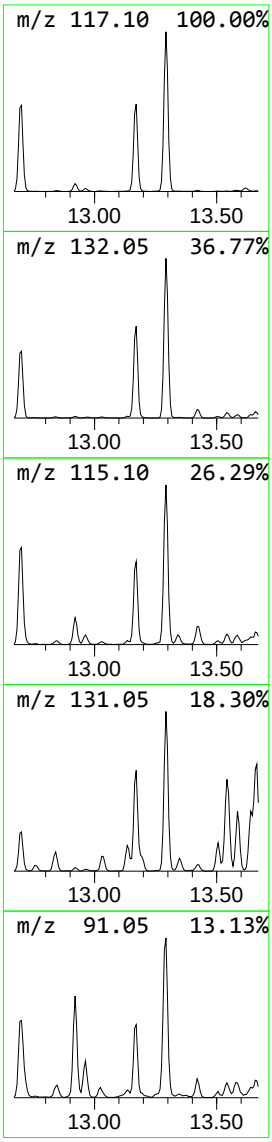
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.170	54.17 ug/l	1187530	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	95
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
5	Indan, 1-methyl-	132	C10H12	000767-58-8	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
Sample : 03019-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

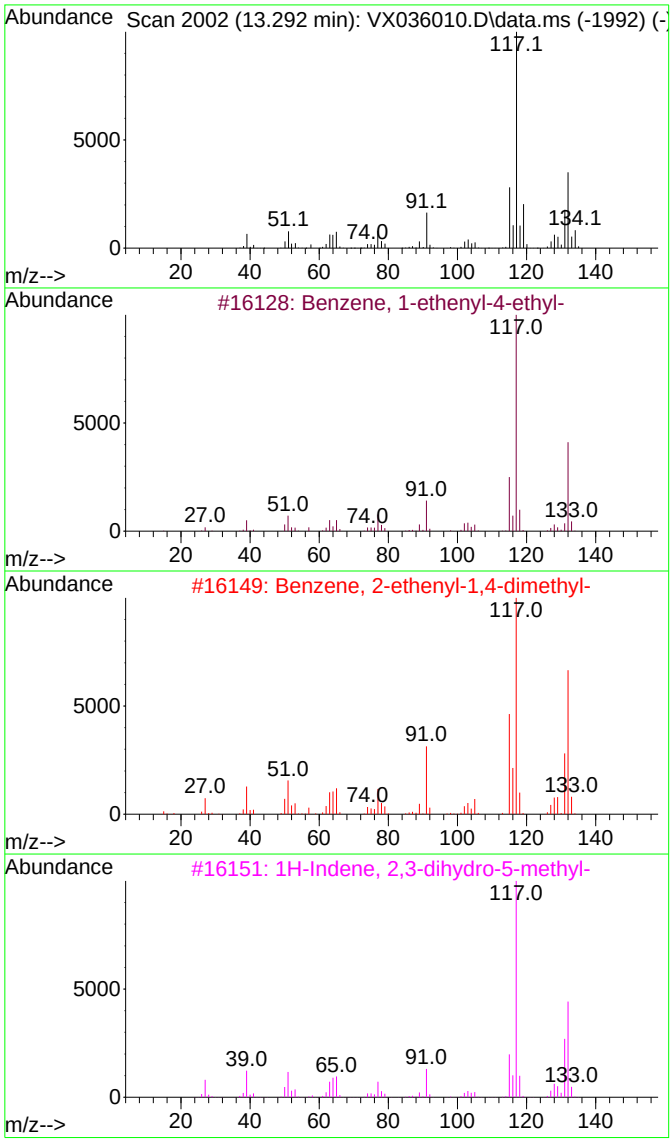
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

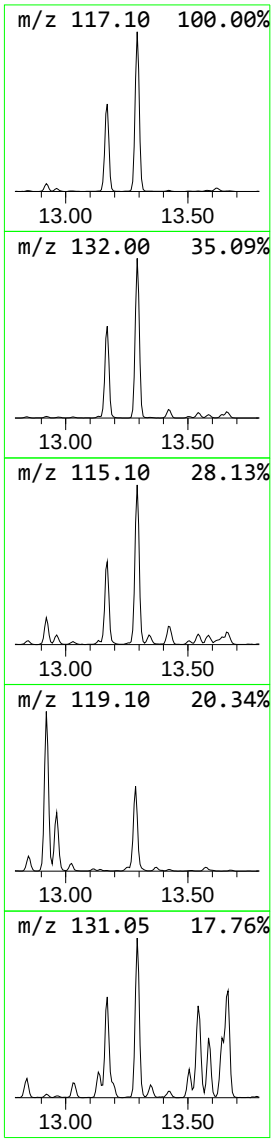
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	120.42 ug/l	2639930	1,4-Dichlorobenzene-d4	12.024

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	91
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5	Indan, 1-methyl-	132	C10H12	000767-58-8	87





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX060523\
Data File : VX036010.D
Acq On : 05 Jun 2023 18:03
Operator : JC/MD
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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MH-289

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X053123W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Ethane, 1,1-dif...	1.142	50.1	ug/l	837720	1	5.556	836005	50.0
Butane, 2-methyl-	1.751	160.1	ug/l	2676870	1	5.556	836005	50.0
Pentane, 2-methyl-	2.824	58.8	ug/l	982821	1	5.556	836005	50.0
Pentane, 3-methyl-	3.099	59.4	ug/l	992538	1	5.556	836005	50.0
Cyclopentane, m...	4.300	98.2	ug/l	1641990	1	5.556	836005	50.0
Indane	12.243	154.6	ug/l	3388460	4	12.024	1096170	50.0
Indan, 1-methyl-	12.701	57.2	ug/l	1253040	4	12.024	1096170	50.0
Benzene, 1,2,3,...	12.920	51.2	ug/l	1122190	4	12.024	1096170	50.0
1H-Indene, 2,3-...	13.170	54.2	ug/l	1187530	4	12.024	1096170	50.0
Benzene, 1-ethe...	13.292	120.4	ug/l	2639930	4	12.024	1096170	50.0