LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

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\82X022825W.M
Title : SW846 8260

Signal : TIC: VX045096.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.740	102	107	123	rVB	304295	416070	10.44%	1.280%
2	1.947	136	141	151	rVB	188563	259099	6.50%	0.797%
3	2.331	195	204	216	rBV	60337	123829	3.11%	0.381%
4	2.776	269	277	280	rBV	159857	351123	8.81%	1.080%
5	2.819	280	284	295	rVB	390685	779975	19.57%	2.399%
6	3.093	317	329	344	rBV	398047	896924	22.51%	2.758%
7	3.440	377	386	396	rBV2	82736	189663	4.76%	0.583%
8	4.044	474	485	490	rBV3	14292	44201	1.11%	0.136%
9	4.209	501	512	518	rBV2	40995	120880	3.03%	0.372%
10	4.300	518	527	539	rVB	291269	833904	20.93%	2.565%
11	5.123	649	662	682	rVB6	19960	96655	2.43%	0.297%
12	5.385	688	705	710	rBV	56409	172865	4.34%	0.532%
13	5.471	711	719	727	rBV3	82148	307184	7.71%	0.945%
14	5.611	737	742	761	rVB2	117762	393322	9.87%	1.210%
15	5.818	765	776	790	rBV3	100674	320585	8.05%	0.986%
16	5.952	790	798	810	rBV	57473	143580	3.60%	0.442%
17	6.153	817	831	840	rBV3	85024	304774	7.65%	0.937%
18	6.251	840	847	856	rVV3	116770	348740	8.75%	1.073%
19	6.355	856	864	875	rVB2	63849	186403	4.68%	0.573%
20	6.605	895	905	918	rBV5	21186	56169	1.41%	0.173%
21	6.757	921	930	937	rBV	127217	336014	8.43%	1.033%
22	6.836	938	943	955	rVV4	67035	210753	5.29%	0.648%
23	7.062	972	980	989	rVB2	18044	42908	1.08%	0.132%
24	7.165	989	997	1005	rBV	40356	91194	2.29%	0.280%
25	7.373	1019	1031	1043	rBV4	194747	531749	13.34%	1.635%
26	7.556	1056	1061	1067	rVV3	28338	60438	1.52%	0.186%
27	7.653	1071	1077	1089	rVB2	44153	93336	2.34%	0.287%
28	7.769	1089	1096	1103	rVB2	39505	78269	1.96%	0.241%
29	7.879	1103	1114	1121	rBV8	16027	56584	1.42%	0.174%
30	7.982	1121	1131	1143	rVB4	46336	145425	3.65%	0.447%
31	8.104	1143	1151	1154	rBV	77064	161989	4.07%	0.498%
32	8.141	1154	1157	1161	rVV	65338	108165	2.71%	0.333%
33	8.184	1161	1164	1175	rVB4	33635	73588	1.85%	0.226%
34	8.312	1175	1185	1191	rBV4	38269	81250	2.04%	0.250%
35	8.507	1212	1217	1226	rVB5	45999	100377	2.52%	0.309%
36	8.598	1226	1232	1235	rBV2	58569	111706	2.80%	0.344%

LSC Area Percent Report

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Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

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\82X022825W.M
Title : SW846 8260

37	8.647	1235	1240	1249	rVB	273499	448105	11.25%	1.378%
38	8.769	1256	1260	1264	rVB2	31797	46130	1.16%	0.142%
39	8.921	1277	1285	1289	rBV2	29334	51618	1.30%	0.159%
40	8.970	1289	1293	1299	rVB	42208	67824	1.70%	0.209%
41	9.098	1306	1314	1319	rBV	58412	104356	2.62%	0.321%
42	9.159	1319	1324	1329	rBV	53685	79582	2.00%	0.245%
43	9.506	1376	1381	1386	rVB4	36093	68141	1.71%	0.210%
44	9.756	1416	1422	1430	rBV5	21042	42796	1.07%	0.132%
45	10.049	1465	1470	1476	rBV	259281	364718	9.15%	1.122%
46	10.299	1508	1511	1517	rVB2	33909	54812	1.38%	0.169%
47	10.640	1563	1567	1581	rVB	29467	49953	1.25%	0.154%
48	10.957	1614	1619	1629	rBV	190072	261052	6.55%	0.803%
49	11.079	1634	1639	1644	rBV	222839	287653	7.22%	0.885%
50	11.299	1668	1675	1683	rBV	791064	1005341	25.23%	3.092%
51	11.396	1683	1691	1696	rVV	1160541	1836744	46.09%	5.649%
52	11.451	1696	1700	1711	rVB	887580	1133854	28.46%	3.487%
53	11.610	1721	1726	1738	rVB	933178	1140374	28.62%	3.507%
54	11.750	1744	1749	1757	rBV	3327533	3984705	100.00%	12.255%
55	11.866	1763	1768	1770	rBV	77450	113978	2.86%	0.351%
56	11.890	1770	1772	1780	rVB	84846	91789	2.30%	0.282%
57	11.969	1780	1785	1788	rBV	52249	63524	1.59%	0.195%
58	12.018	1788	1793	1799	rVV2	297958	400372	10.05%	1.231%
59	12.085	1799	1804	1815	rVB	643337	794694	19.94%	2.444%
60	12.244	1824	1830	1837	rBV	526989	753736	18.92%	2.318%
61	12.311	1837	1841	1851	rVB3	587741	963521	24.18%	2.963%
62	12.402	1851	1856	1861	rVB	90980	110952	2.78%	0.341%
63	12.463	1861	1866	1872	rBV2	321646	420462	10.55%	1.293%
64	12.530	1872	1877	1879	rBV	437565	548822	13.77%	1.688%
65	12.555	1879	1881	1885	rVV	281104	292366	7.34%	0.899%
66	12.610	1885	1890	1894	rVV	1026307	1219947	30.62%	3.752%
67	12.640	1894	1895	1900	rVB	157584	156960	3.94%	0.483%
68	12.695	1900	1904	1912	rBV2	445474	679387	17.05%	2.089%
69	12.847	1924	1929	1934	rVB	195670	245255	6.15%	0.754%
70	12.920	1934	1941	1944	rBV	524636	652564	16.38%	2.007%
71	12.963	1944	1948	1954	rVB	513076	636239	15.97%	1.957%
72	13.024	1954	1958	1964	rBV3	112624	185291	4.65%	0.570%
73	13.164	1978	1981	1986	rVB	398214	468230	11.75%	1.440%
74	13.256	1991	1996	1998	rBV	93060	138335	3.47%	0.425%
75	13.292	1998	2002	2007	rVB2	748054	992395	24.91%	3.052%
76	13.366	2012	2014	2020	rVV	97024	127614	3.20%	0.392%

LSC Area Percent Report

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Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

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\82X022825W.M
Title : SW846 8260

77	13.420	2020	2023	2031	rVB	94772	122660	3.08%	0.377%
78	13.506	2031	2037	2039	rBV2	80074	111506	2.80%	0.343%
79	13.542	2039	2043	2046	rVV	173618	247124	6.20%	0.760%
80	13.579	2046	2049	2056	rVV3	232948	429012	10.77%	1.319%
81	13.664	2056	2063	2069	rVB2	149812	342647	8.60%	1.054%
82	13.792	2079	2084	2090	rVB3	21924	41881	1.05%	0.129%
83	13.926	2099	2106	2110	rBV3	86769	136624	3.43%	0.420%
84	13.969	2110	2113	2117	rVB	63138	71939	1.81%	0.221%
85	14.073	2125	2130	2136	rBV	137657	171442	4.30%	0.527%
86	14.213	2148	2153	2159	rBV	110239	183003	4.59%	0.563%
87	14.317	2167	2170	2175	rBV	54511	70281	1.76%	0.216%
88	14.371	2175	2179	2183	rVB	77350	93604	2.35%	0.288%
89	14.634	2209	2222	2233	rBV	384780	503392	12.63%	1.548%
90	14.780	2240	2246	2256	rVB	207324	276421	6.94%	0.850%

Sum of corrected areas: 32515417

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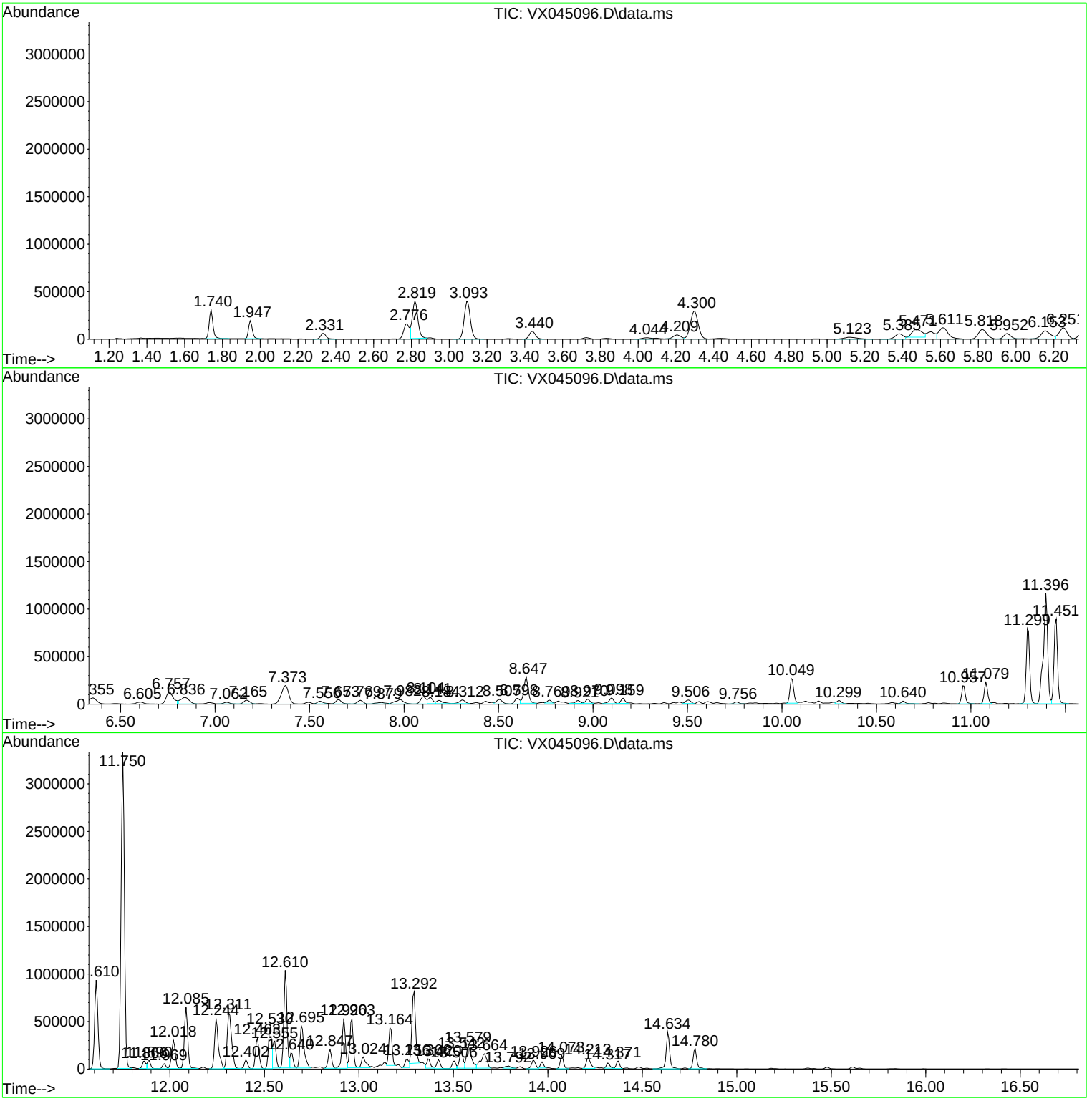
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

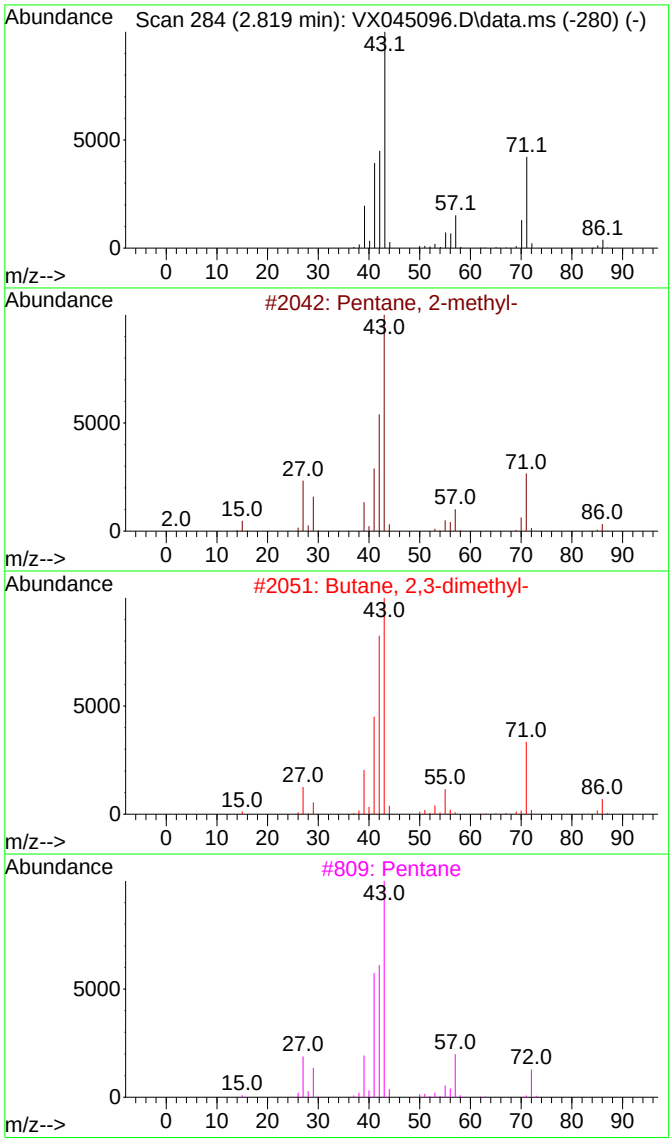
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

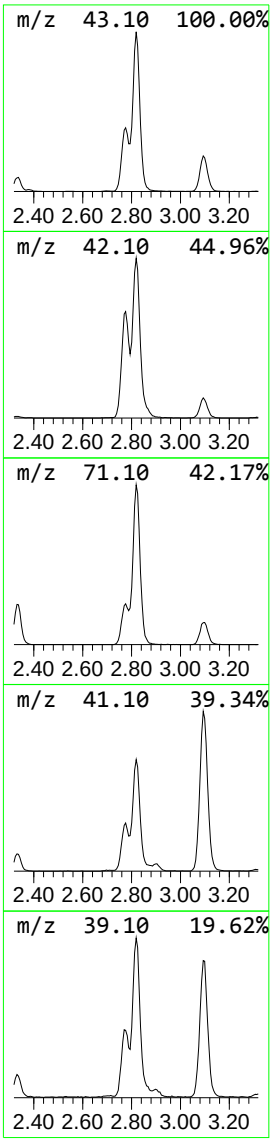
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Pentane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.819	99.15 ug/l	779975	Pentafluorobenzene	5.550

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	46
4	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45
5	Heptane, 4-methyl-	114	C8H18	000589-53-7	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

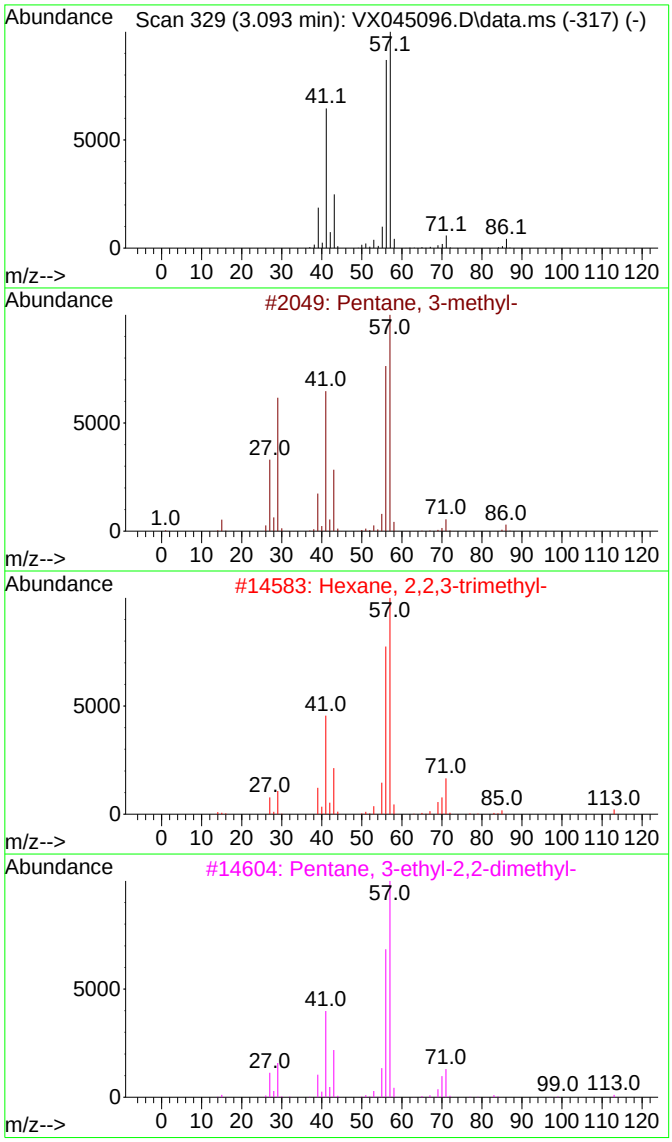
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

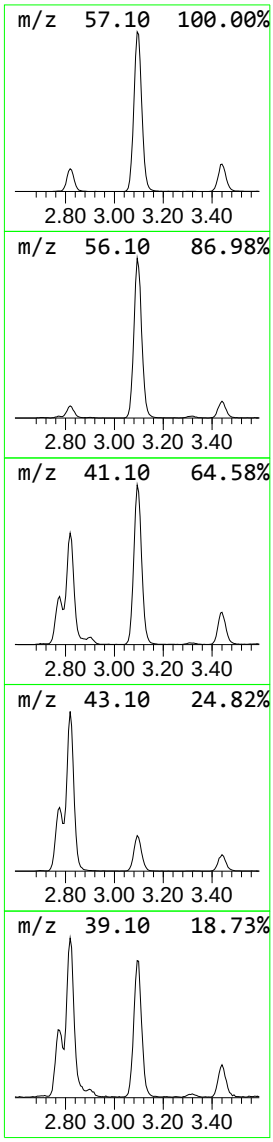
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.093	114.02 ug/l	896924	Pentafluorobenzene	5.550

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	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64
4	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	50
5	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

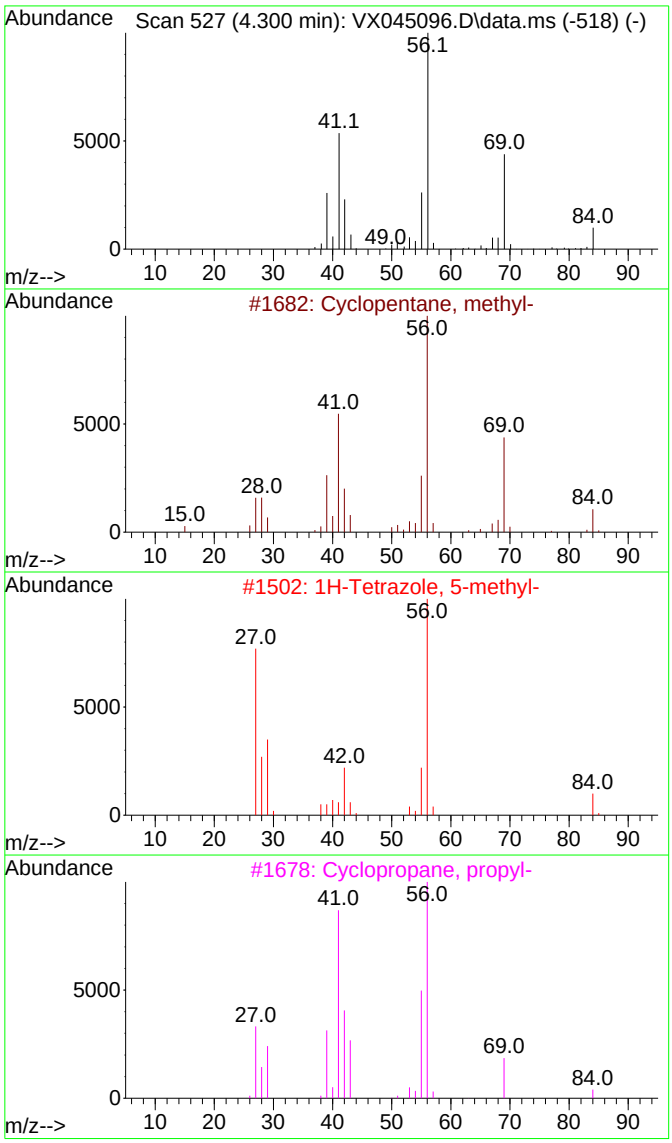
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

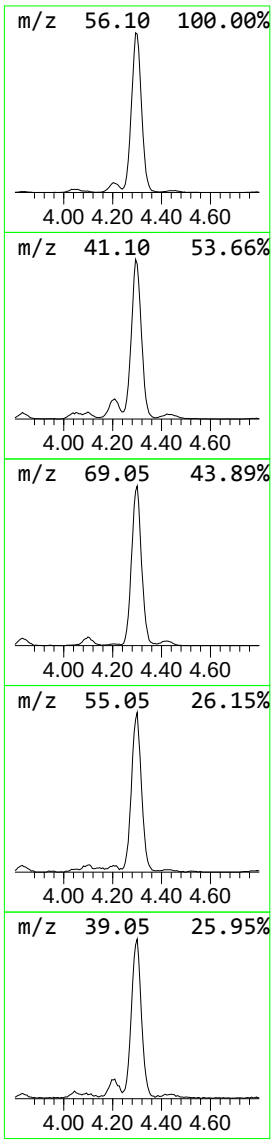
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.300	106.01 ug/l	833904	Pentafluorobenzene	5.550

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	56
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	56
4	Cyclobutane	56	C4H8	000287-23-0	50
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

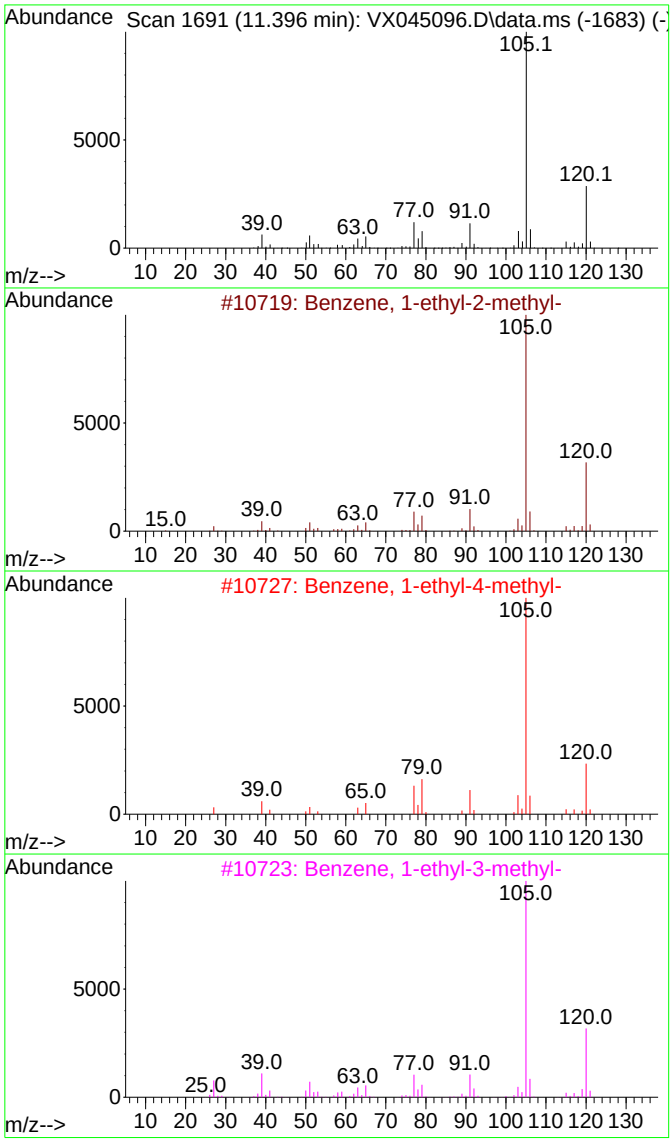
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

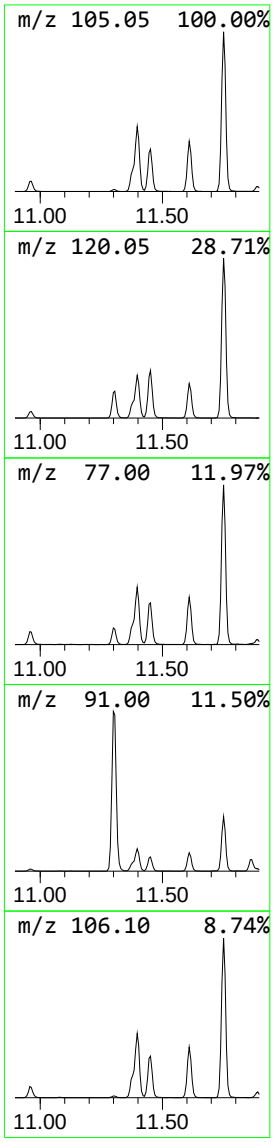
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.396	229.38 ug/l	1836740	1,4-Dichlorobenzene-d4	12.018

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-4-methyl-	120	C9H12	000622-96-8	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

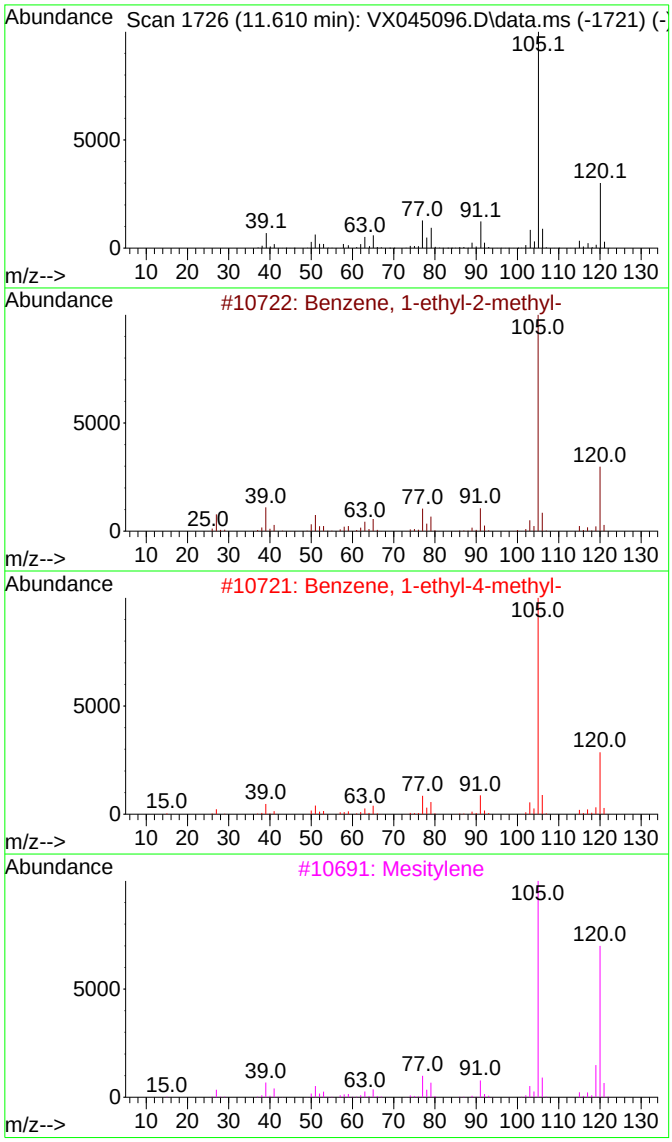
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

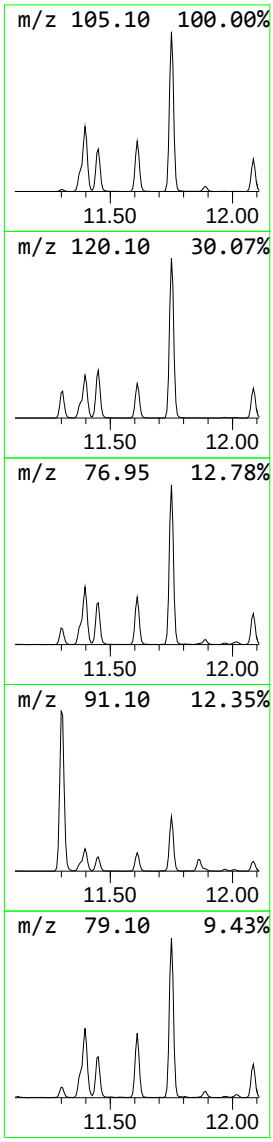
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-4-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	142.41 ug/l	1140370	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

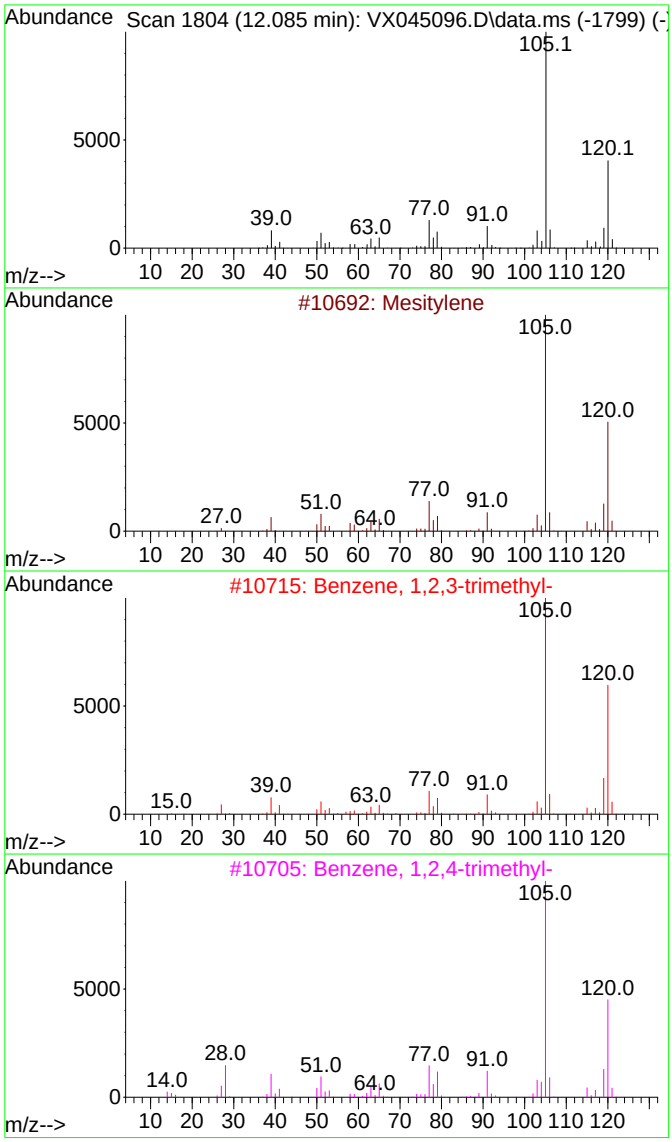
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

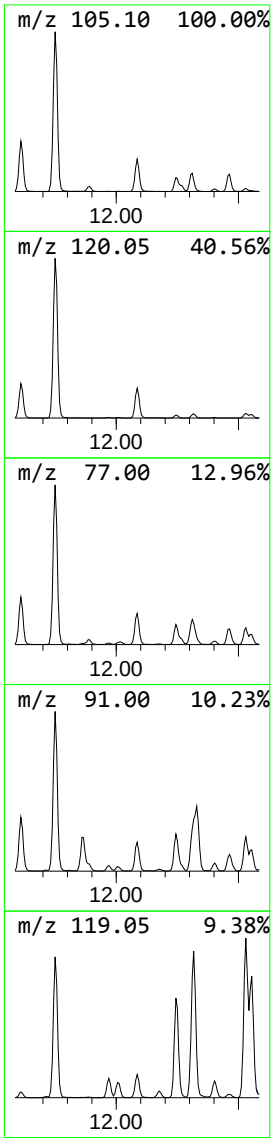
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1,2,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.085	99.24 ug/l	794694	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

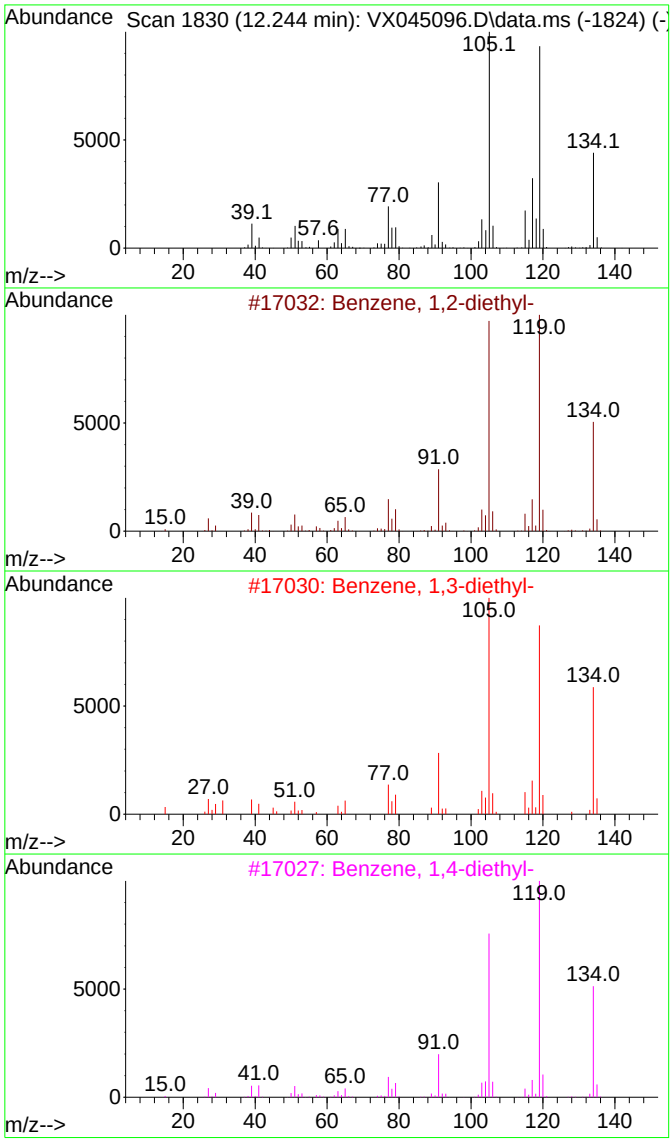
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

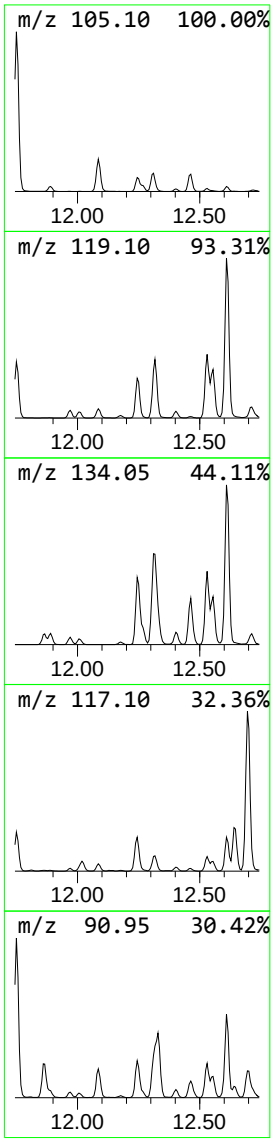
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,2-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	94.13 ug/l	753736	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	90
3	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	68
5	Ethanone, 1-(4-methylphenyl)-	134	C9H10O	000122-00-9	60





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

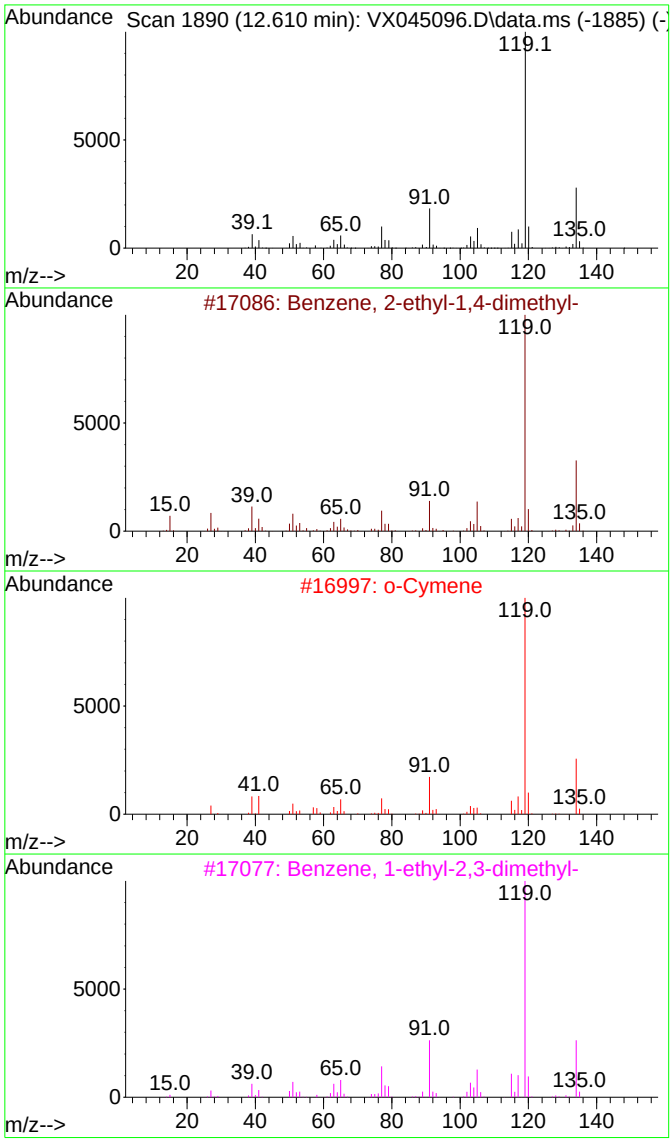
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

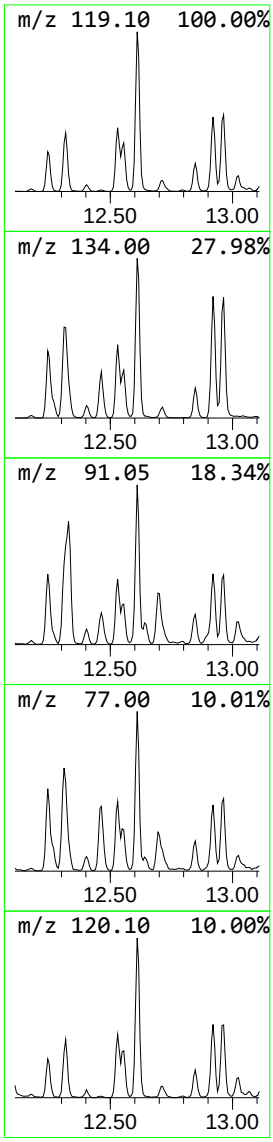
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.610	152.35 ug/l	1219950	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

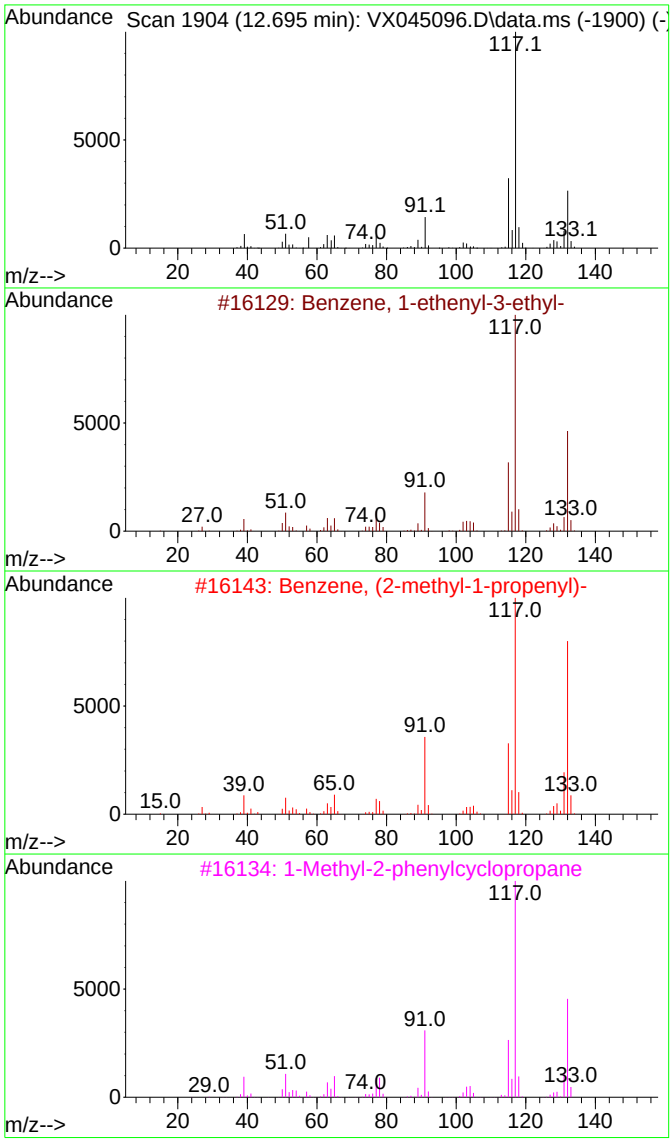
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

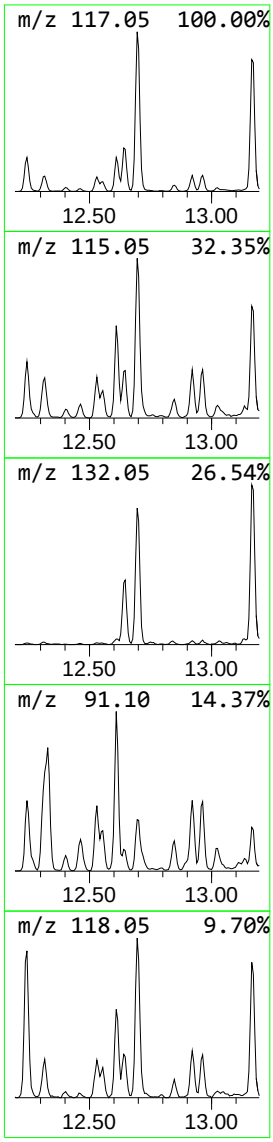
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	84.84 ug/l	679387	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	90
2	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	90
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	87
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

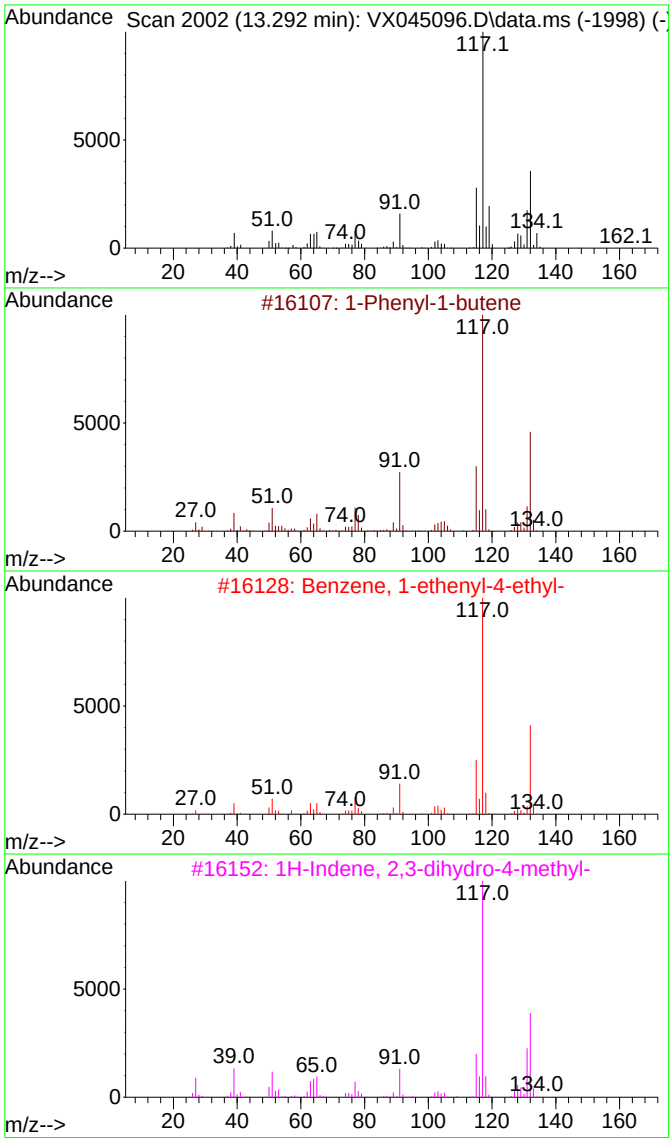
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
Quant Title : SW846 8260

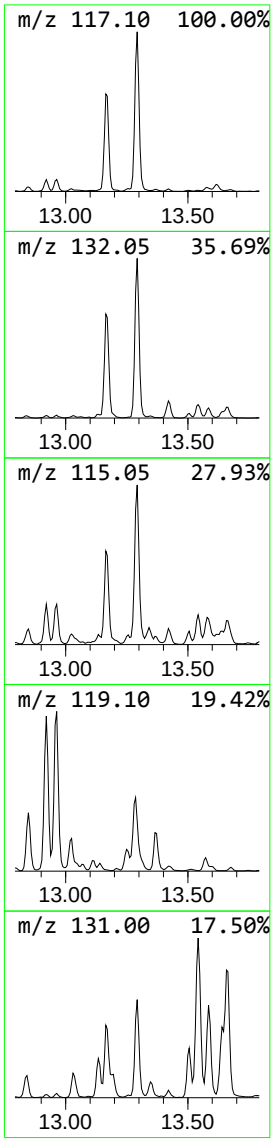
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.292	123.93 ug/l	992395	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	89
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	89
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	81
5	Indan, 1-methyl-	132	C10H12	000767-58-8	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022825\
Data File : VX045096.D
Acq On : 28 Feb 2025 18:04
Operator : JC/MD
Sample : Q1462-01
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
MW2

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.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
Pentane, 2-methyl-	2.819	99.2	ug/l	779975	1	5.550	393322	50.0
Pentane, 3-methyl-	3.093	114.0	ug/l	896924	1	5.550	393322	50.0
Cyclopentane, m...	4.300	106.0	ug/l	833904	1	5.550	393322	50.0
Benzene, 1-ethy...	11.396	229.4	ug/l	1836740	4	12.018	400372	50.0
Benzene, 1-ethy...	11.610	142.4	ug/l	1140370	4	12.018	400372	50.0
Benzene, 1,2,3-...	12.085	99.2	ug/l	794694	4	12.018	400372	50.0
Benzene, 1,2-di...	12.244	94.1	ug/l	753736	4	12.018	400372	50.0
Benzene, 2-ethy...	12.610	152.3	ug/l	1219950	4	12.018	400372	50.0
Benzene, 1-ethe...	12.695	84.8	ug/l	679387	4	12.018	400372	50.0
1-Phenyl-1-butene	13.292	123.9	ug/l	992395	4	12.018	400372	50.0