LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

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\82X070225W.M
Title : SW846 8260

Signal : TIC: VX046944.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.264	25	29	42	rVB3	11480	29544	1.08%	0.114%
2	1.770	105	112	128	rBV	212265	349338	12.74%	1.344%
3	2.361	198	209	220	rBV	81937	188371	6.87%	0.725%
4	2.800	272	281	292	rBV	238346	622395	22.69%	2.395%
5	3.117	324	333	350	rBV	167137	418139	15.24%	1.609%
6	4.062	479	488	500	rBV4	11644	41138	1.50%	0.158%
7	4.227	500	515	525	rVV5	20823	75814	2.76%	0.292%
8	4.446	539	551	564	rBV7	21136	83664	3.05%	0.322%
9	5.147	654	666	681	rBV9	21979	103909	3.79%	0.400%
10	5.397	694	707	723	rBV2	270661	855406	31.19%	3.291%
11	5.562	724	734	743	rBV	424877	1335388	48.68%	5.138%
12	5.861	772	783	792	rBV3	47722	156343	5.70%	0.602%
13	5.964	792	800	816	rVV	288762	802329	29.25%	3.087%
14	6.196	829	838	843	rBV7	14307	48672	1.77%	0.187%
15	6.269	843	850	858	rVV3	29652	87792	3.20%	0.338%
16	6.367	858	866	880	rVB3	62644	188189	6.86%	0.724%
17	6.769	923	932	960	rBV	746324	1971795	71.89%	7.587%
18	6.982	960	967	976	rVB5	18602	49261	1.80%	0.190%
19	7.178	991	999	1010	rBV3	61657	151986	5.54%	0.585%
20	7.367	1021	1030	1041	rVB3	58317	153846	5.61%	0.592%
21	7.507	1044	1053	1058	rBV	21747	48945	1.78%	0.188%
22	7.562	1058	1062	1074	rVB3	23161	52230	1.90%	0.201%
23	7.781	1090	1098	1106	rVB3	38923	85909	3.13%	0.331%
24	8.007	1123	1135	1144	rVB4	23417	86616	3.16%	0.333%
25	8.110	1144	1152	1161	rBV3	40197	90462	3.30%	0.348%
26	8.220	1161	1170	1179	rVB6	16686	43453	1.58%	0.167%
27	8.311	1179	1185	1194	rBV5	48809	119326	4.35%	0.459%
28	8.513	1209	1218	1227	rVB5	32812	82580	3.01%	0.318%
29	8.647	1227	1240	1253	rBV	1593828	2742969	100.00%	10.554%
30	8.775	1255	1261	1265	rBV	29733	47718	1.74%	0.184%
31	8.921	1281	1285	1289	rVB2	20091	30517	1.11%	0.117%
32	8.976	1289	1294	1300	rVB2	66782	111582	4.07%	0.429%
33	9.104	1307	1315	1320	rVV4	61080	119310	4.35%	0.459%
34	9.165	1320	1325	1335	rVB	208205	355664	12.97%	1.368%
35	9.323	1344	1351	1357	rBV5	14558	32679	1.19%	0.126%
36	9.427	1363	1368	1371	rBV	110809	174830	6.37%	0.673%

LSC Area Percent Report

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Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

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\82X070225W.M
Title : SW846 8260

37	9.458	1371	1373	1376	rVV	79170	108445	3.95%	0.417%
38	9.500	1376	1380	1389	rVB3	95354	184335	6.72%	0.709%
39	9.659	1400	1406	1415	rVB4	23140	52677	1.92%	0.203%
40	9.762	1417	1423	1431	rBV2	72730	124158	4.53%	0.478%
41	9.939	1444	1452	1455	rBV3	21887	45744	1.67%	0.176%
42	9.976	1455	1458	1462	rVV4	22301	28440	1.04%	0.109%
43	10.055	1465	1471	1477	rVV	1659214	2288244	83.42%	8.804%
44	10.128	1480	1483	1485	rVV	52023	70198	2.56%	0.270%
45	10.159	1485	1488	1493	rVB2	49642	63192	2.30%	0.243%
46	10.275	1502	1507	1517	rVB5	44529	101424	3.70%	0.390%
47	10.457	1532	1537	1547	rVB3	19159	41544	1.51%	0.160%
48	10.585	1547	1558	1565	rBV4	29219	79611	2.90%	0.306%
49	10.774	1583	1589	1592	rBV3	24672	39619	1.44%	0.152%
50	10.866	1598	1604	1610	rBV7	15441	30926	1.13%	0.119%
51	10.963	1615	1620	1624	rBV	169175	235552	8.59%	0.906%
52	11.079	1634	1639	1644	rBV	1410543	1785701	65.10%	6.871%
53	11.122	1644	1646	1653	rVB4	39699	53379	1.95%	0.205%
54	11.305	1671	1676	1685	rVB	216553	280073	10.21%	1.078%
55	11.616	1722	1727	1734	rVB2	27670	44480	1.62%	0.171%
56	11.713	1738	1743	1745	rBV2	24203	29817	1.09%	0.115%
57	11.750	1745	1749	1754	rVB	48690	67080	2.45%	0.258%
58	11.860	1762	1767	1770	rBV	160453	229313	8.36%	0.882%
59	11.890	1770	1772	1777	rVB	117672	125623	4.58%	0.483%
60	12.018	1788	1793	1802	rBV	1935166	2332730	85.04%	8.975%
61	12.177	1815	1819	1825	rVB	77099	98597	3.59%	0.379%
62	12.244	1825	1830	1837	rBV	583891	719008	26.21%	2.766%
63	12.311	1837	1841	1849	rVB	465105	604972	22.06%	2.328%
64	12.402	1851	1856	1863	rBV	190031	230680	8.41%	0.888%
65	12.530	1871	1877	1879	rBV2	24218	33362	1.22%	0.128%
66	12.609	1884	1890	1894	rVV2	47229	76231	2.78%	0.293%
67	12.658	1894	1898	1900	rVV3	23924	36235	1.32%	0.139%
68	12.695	1900	1904	1911	rVV	346436	499998	18.23%	1.924%
69	12.756	1911	1914	1919	rVB	55729	67834	2.47%	0.261%
70	12.835	1919	1927	1933	rVB2	130290	204503	7.46%	0.787%
71	12.920	1933	1941	1945	rBV	652559	873905	31.86%	3.362%
72	12.963	1945	1948	1954	rVV3	49171	83586	3.05%	0.322%
73	13.030	1954	1959	1964	rVB3	138938	216502	7.89%	0.833%
74	13.134	1966	1976	1980	rBV2	218301	345077	12.58%	1.328%
75	13.189	1982	1985	1991	rVB	127141	167442	6.10%	0.644%
76	13.292	1998	2002	2006	rBV2	41703	55166	2.01%	0.212%

LSC Area Percent Report

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Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

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\82X070225W.M
Title : SW846 8260

77	13.341	2006	2010	2018	rVV3	78927	127688	4.66%	0.491%
78	13.426	2020	2024	2029	rVB4	29544	44504	1.62%	0.171%
79	13.512	2033	2038	2040	rBV5	18862	30808	1.12%	0.119%
80	13.542	2040	2043	2046	rVV	94355	132448	4.83%	0.510%
81	13.573	2046	2048	2052	rVV	112803	146609	5.34%	0.564%
82	13.615	2052	2055	2061	rVV	102226	148839	5.43%	0.573%
83	13.750	2071	2077	2080	rBV2	28485	38984	1.42%	0.150%
84	13.792	2080	2084	2090	rVV4	23345	50758	1.85%	0.195%
85	13.853	2090	2094	2098	rVB	58618	79461	2.90%	0.306%
86	13.926	2098	2106	2110	rBV2	65120	105178	3.83%	0.405%
87	13.969	2110	2113	2118	rVB2	26725	33077	1.21%	0.127%
88	14.219	2149	2154	2162	rBV3	32160	66218	2.41%	0.255%
89	14.316	2166	2170	2177	rVV2	87256	143048	5.22%	0.550%
90	14.597	2210	2216	2220	rBV2	47764	78216	2.85%	0.301%
91	14.780	2241	2246	2256	rVB3	42243	71490	2.61%	0.275%

Sum of corrected areas: 25990838

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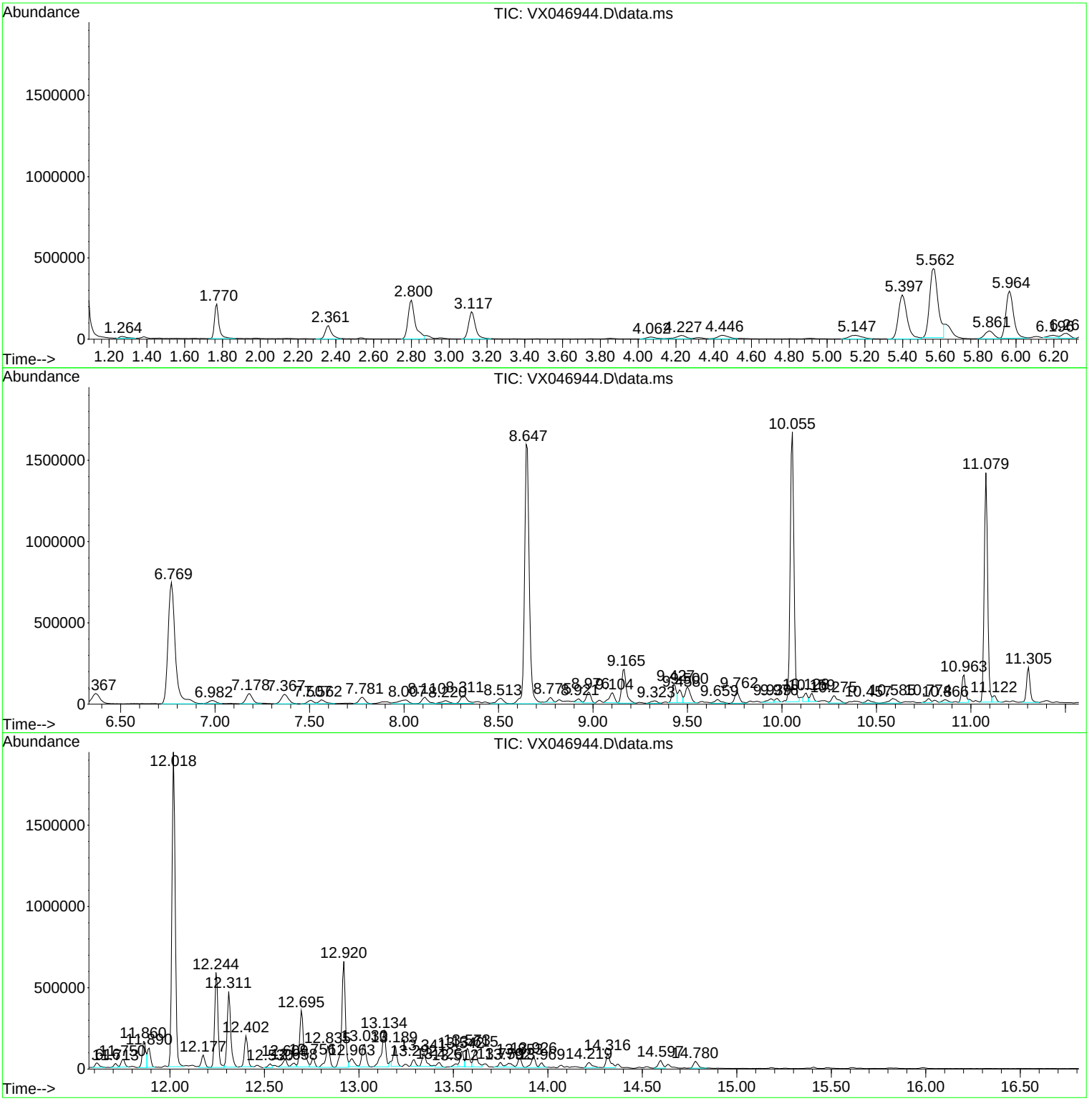
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
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Instrument :
MSVOA_X
ClientSampleId :
AOC-202

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

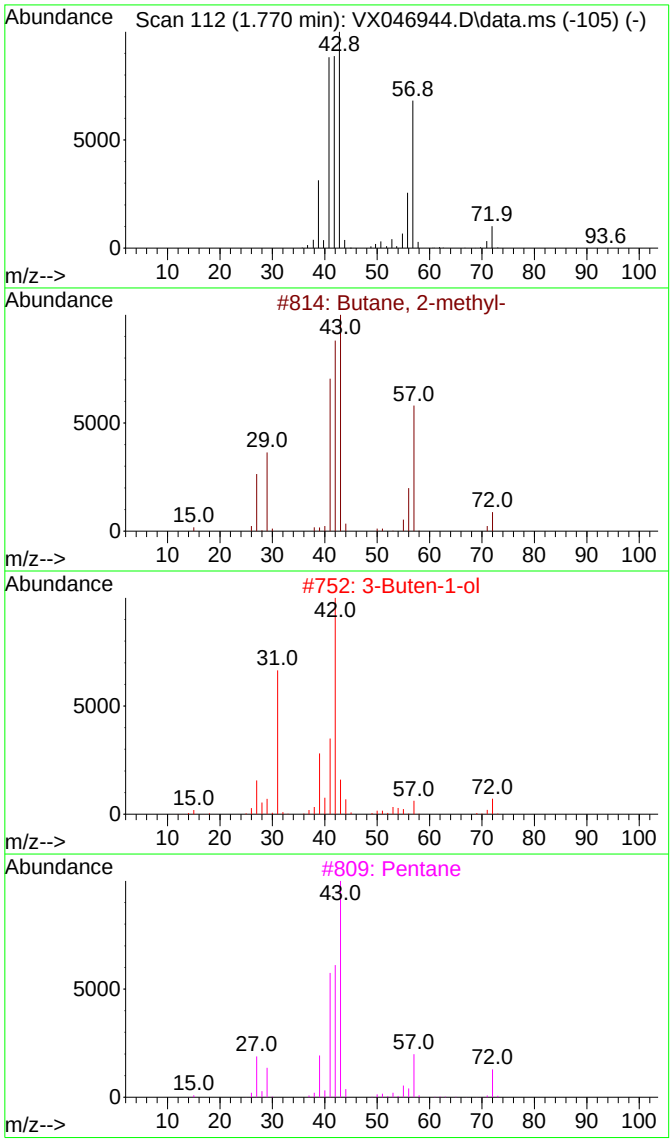
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

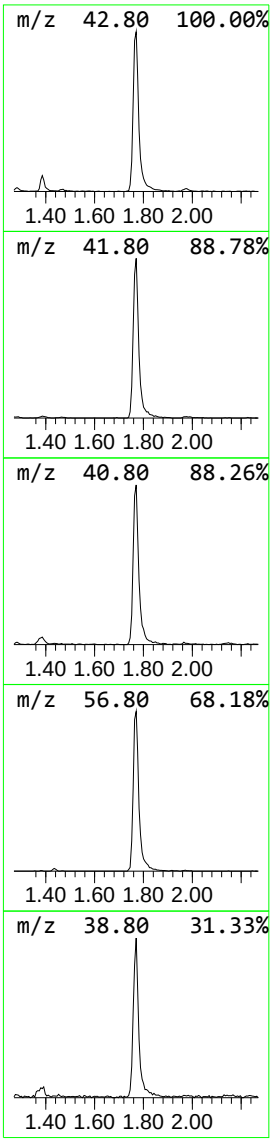
****

Peak Number 1 Butane, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.770	13.08 ug/l	349338	Pentafluorobenzene	5.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	33
4	2,5-Furandione, dihydro-3-methyl-	114	C5H6O3	004100-80-5	23
5	1-Butene	56	C4H8	000106-98-9	18





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

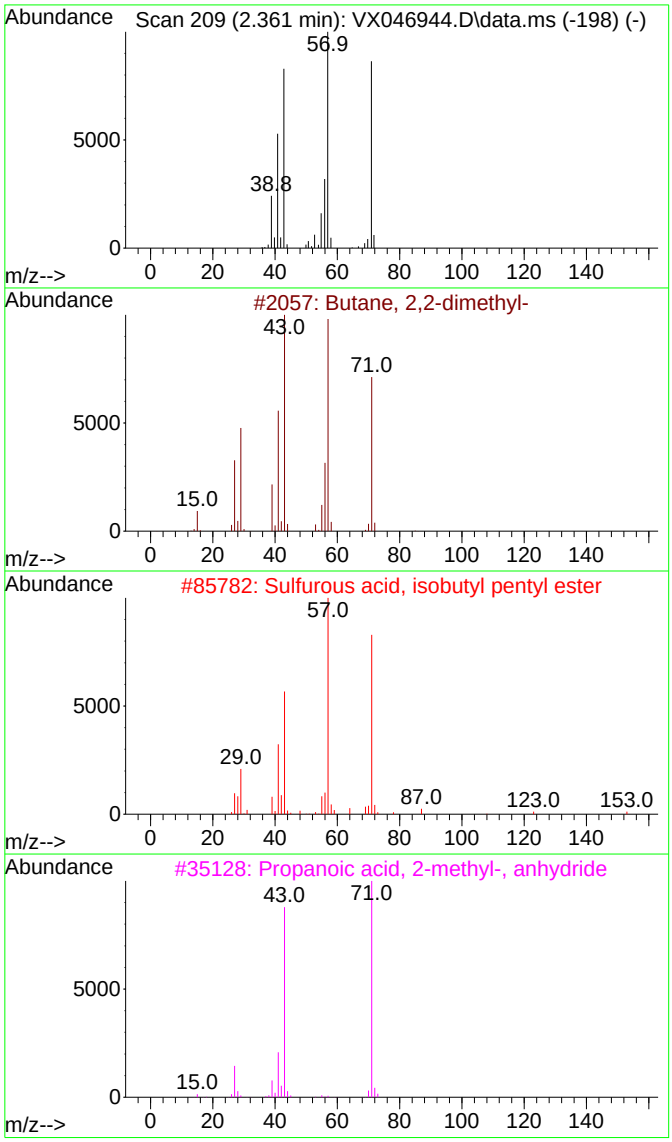
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

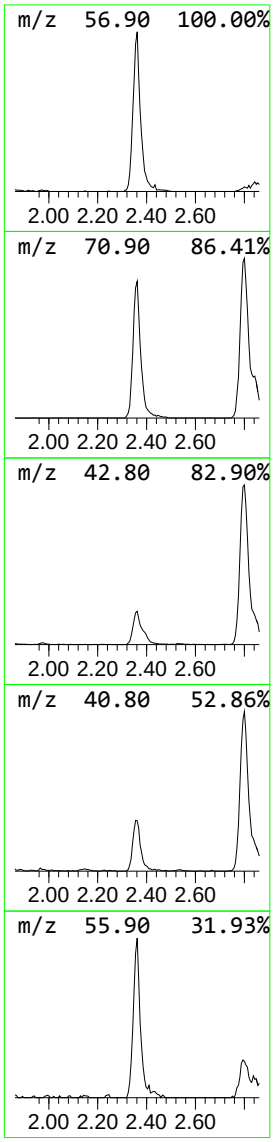
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.361	7.05 ug/l	188371	Pentafluorobenzene	5.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	83
2	Sulfurous acid, isobutyl pentyl ...	208	C9H20O3S	1000309-13-8	64
3	Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	40
4	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	33
5	Oxirane, propyl-	86	C5H10O	001003-14-1	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

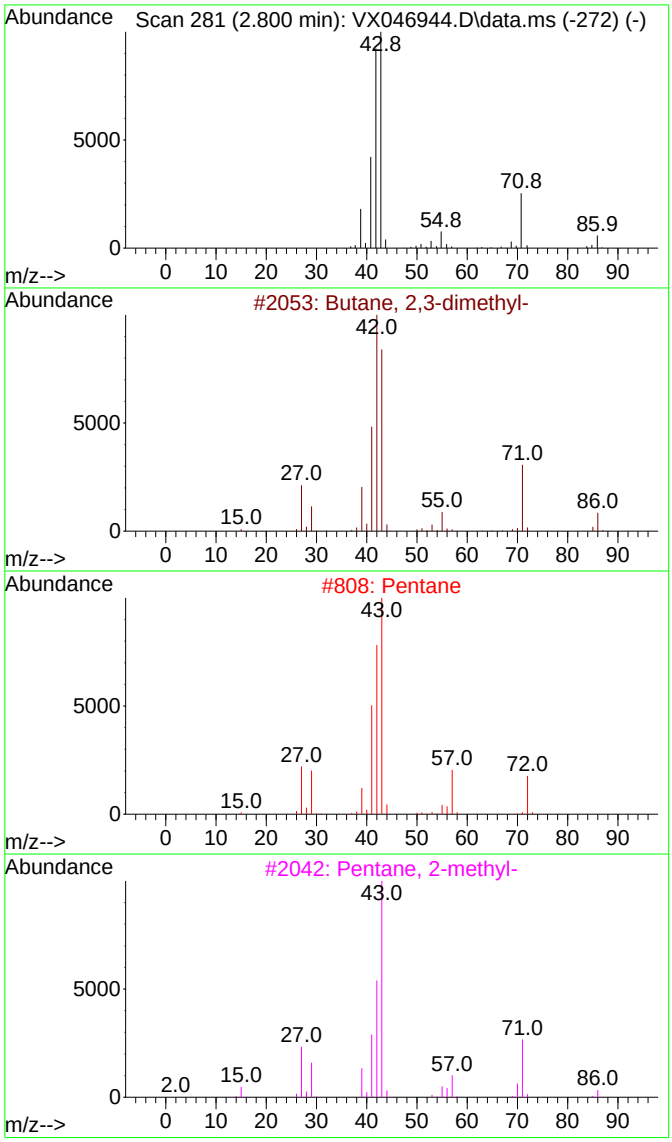
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

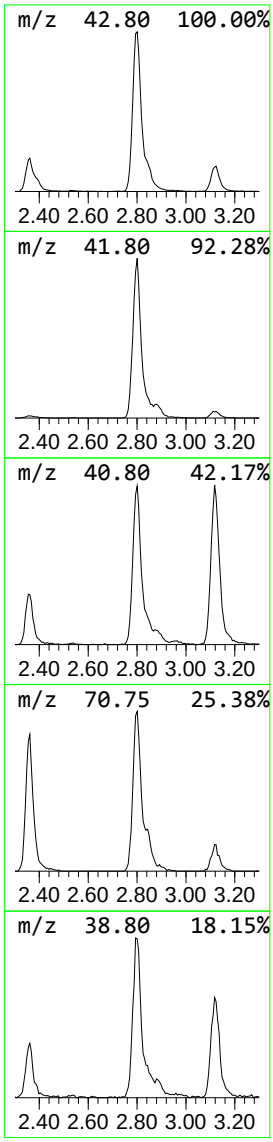
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Butane, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.800	23.30 ug/l	622395	Pentafluorobenzene	5.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	87
2	Pentane	72	C5H12	000109-66-0	45
3	Pentane, 2-methyl-	86	C6H14	000107-83-5	40
4	Tetrahydrofuran	72	C4H8O	000109-99-9	38
5	Propanoyl chloride, 2-methyl-	106	C4H7ClO	000079-30-1	27





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
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Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

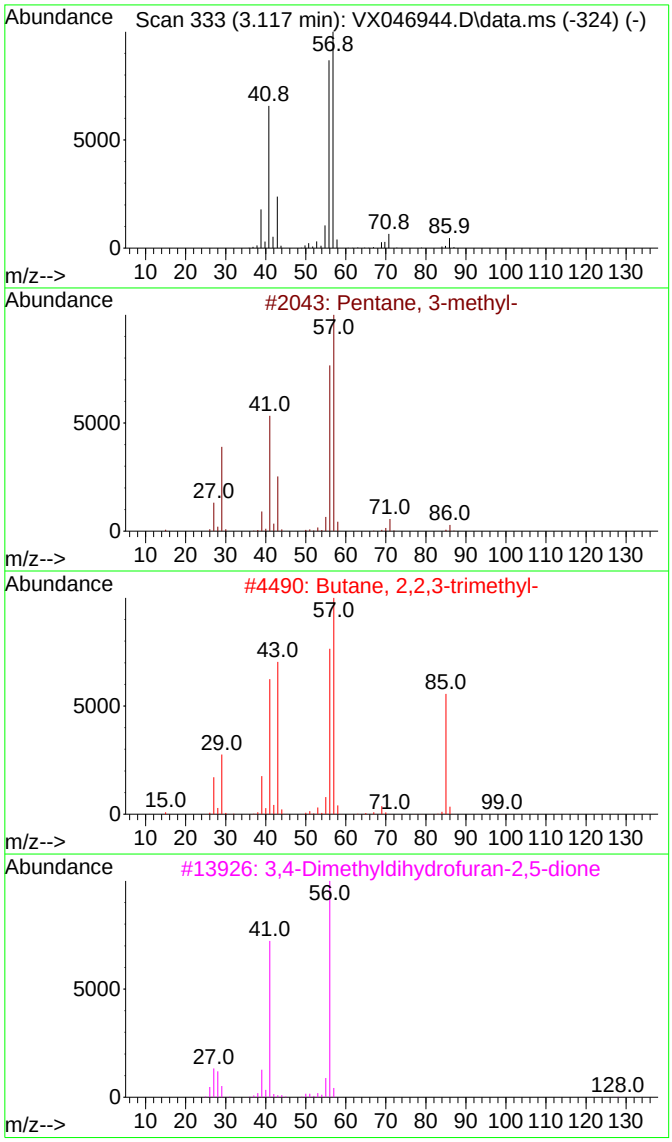
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

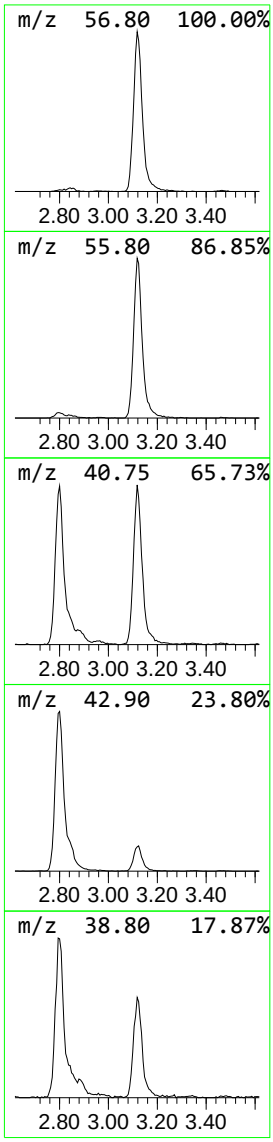
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.117	15.66 ug/l	418139	Pentafluorobenzene	5.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
4	Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	50
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

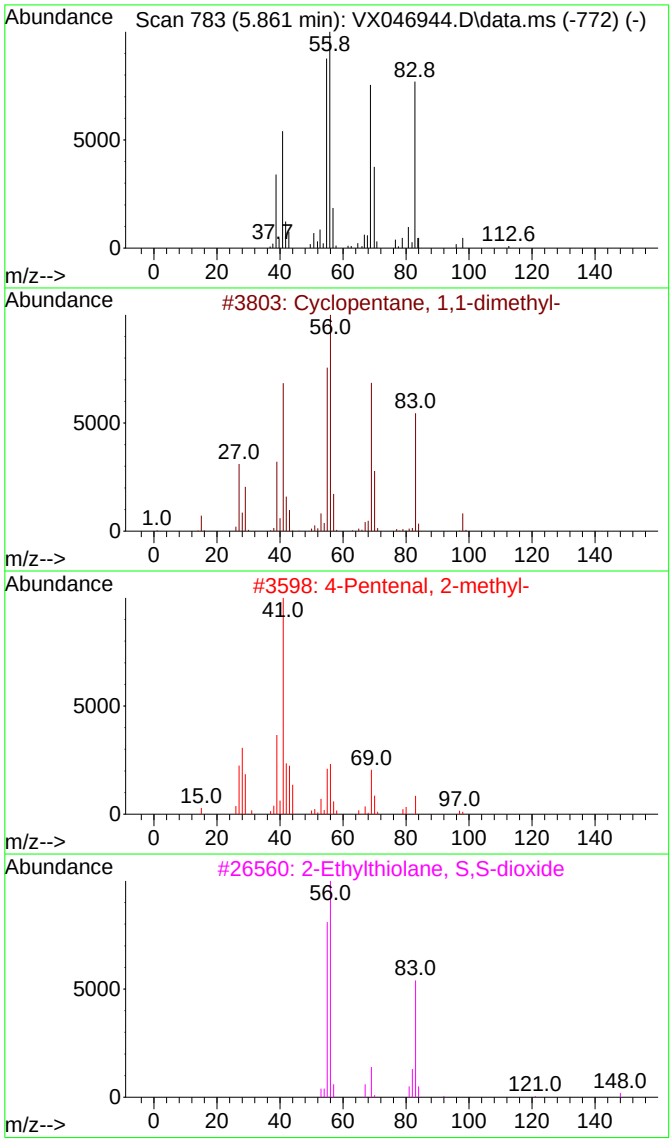
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

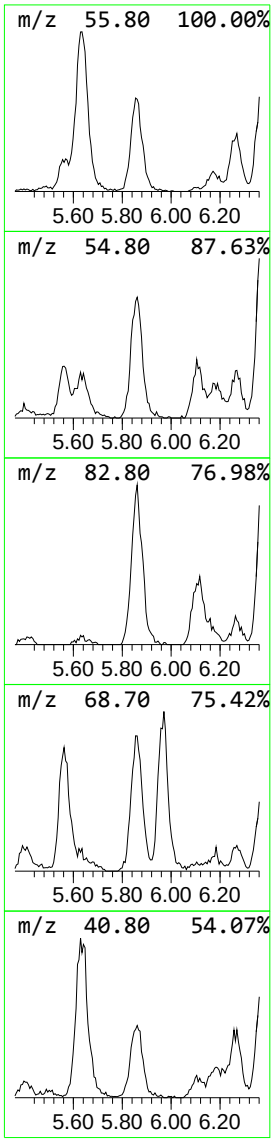
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, 1,1-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.861	5.85 ug/l	156343	Pentafluorobenzene	5.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	78
2	4-Pentenal, 2-methyl-	98	C6H10O	005187-71-3	53
3	2-Ethylthiolane, S,S-dioxide	148	C6H12O2S	010178-59-3	50
4	2-Heptene	98	C7H14	000592-77-8	50
5	1-Heptene, 6-methyl-	112	C8H16	005026-76-6	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

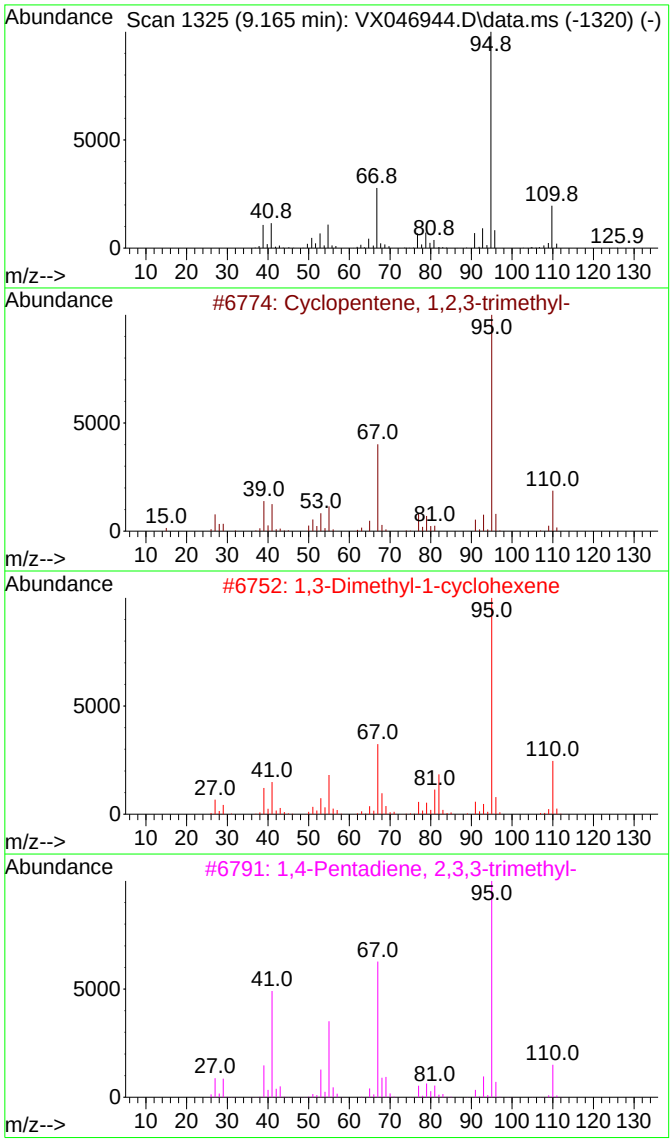
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

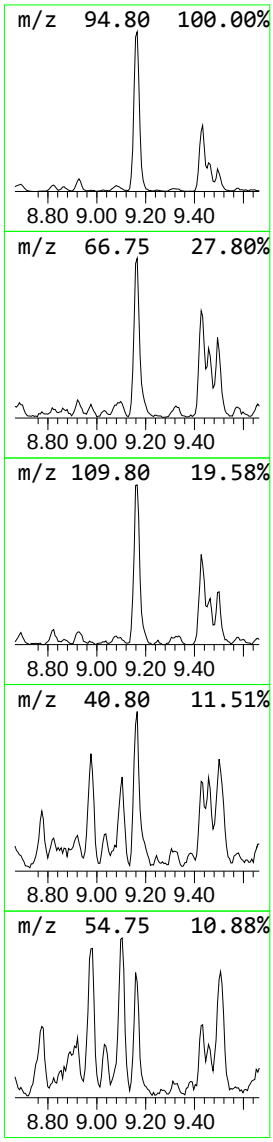
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.165	7.77 ug/l	355664	Chlorobenzene-d5	10.055

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	91
2	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	91
3	1,4-Pentadiene, 2,3,3-trimethyl-	110	C8H14	000756-02-5	78
4	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1010142-17-5	72
5	1,4-Pentadiene, 2,3,4-trimethyl-	110	C8H14	072014-90-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

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

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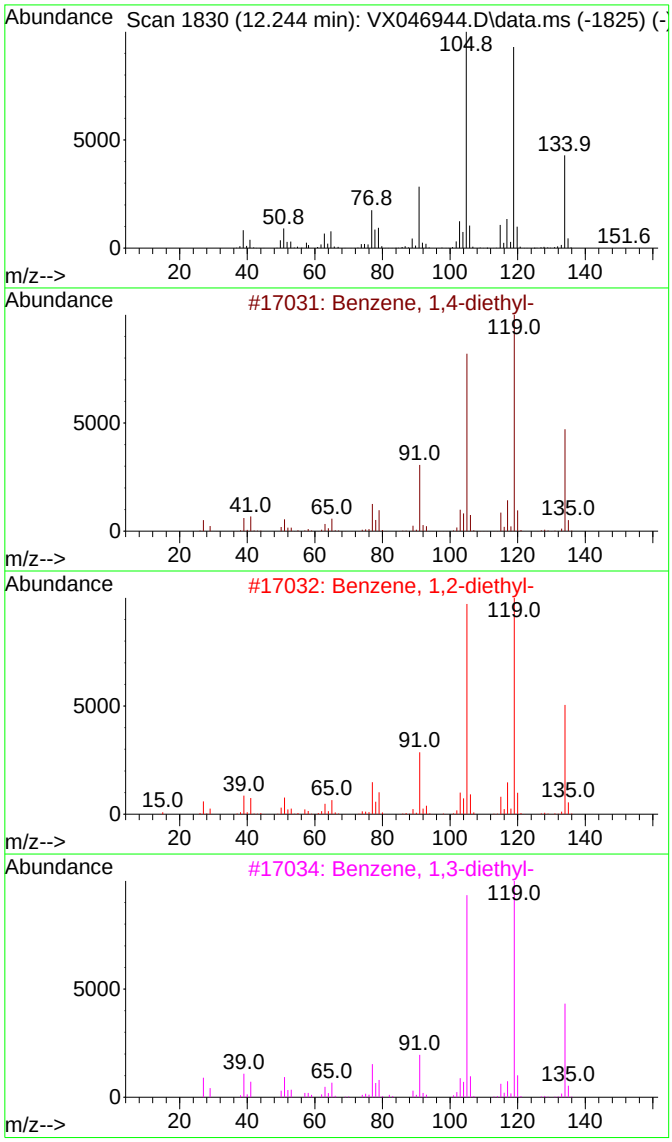
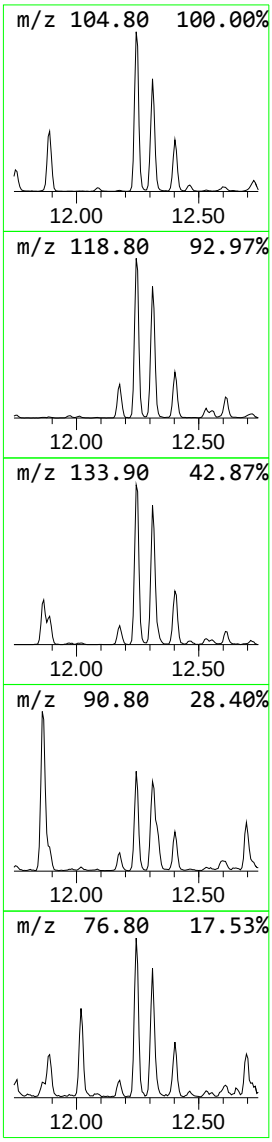
Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.244	15.41 ug/l	719008	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
5	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

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

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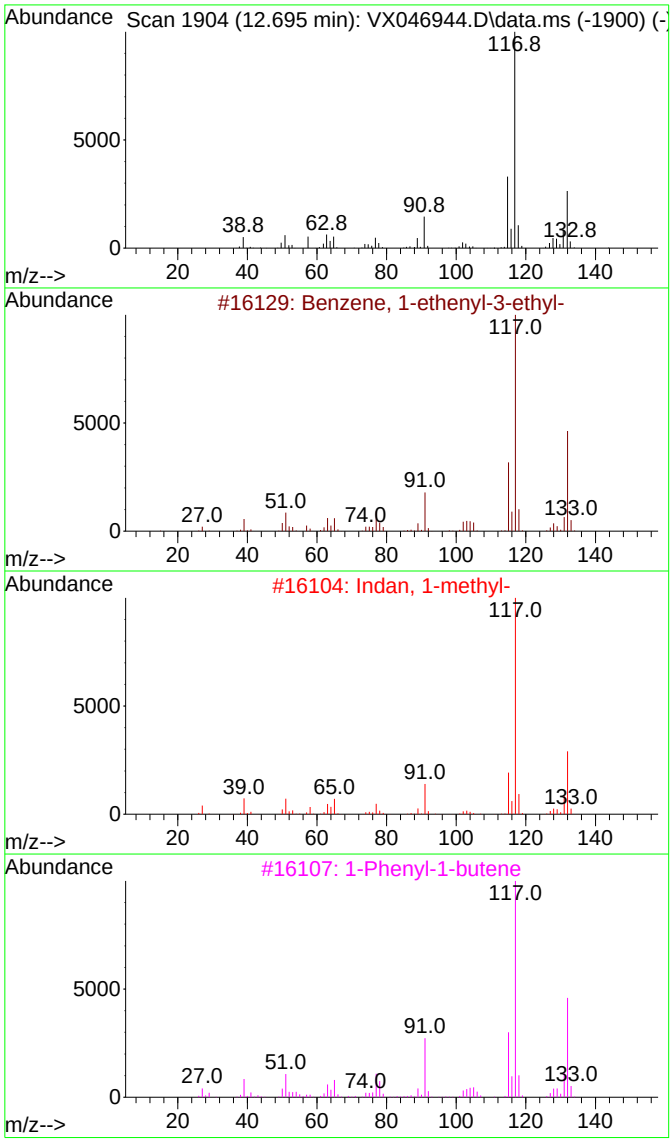
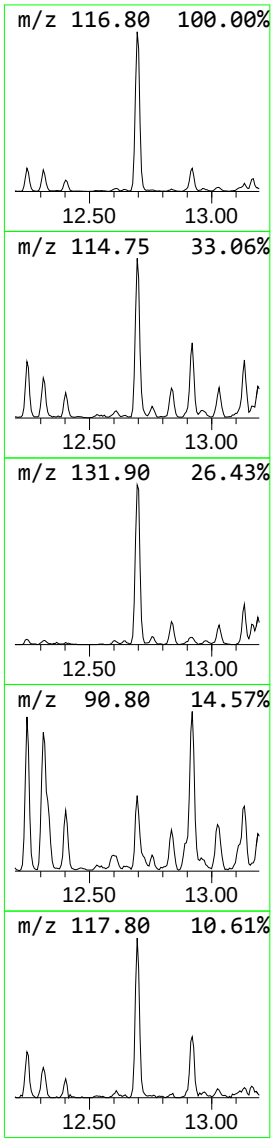
Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.695	10.72 ug/l	499998	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	91
2	Indan, 1-methyl-	132	C10H12	000767-58-8	90
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	90
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

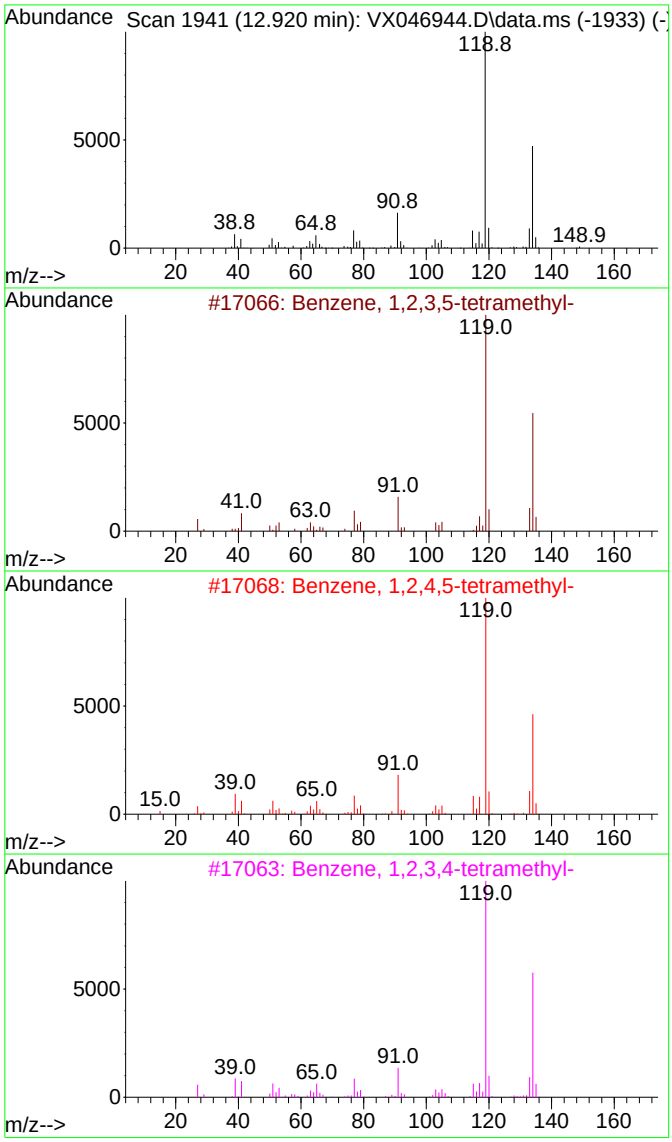
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

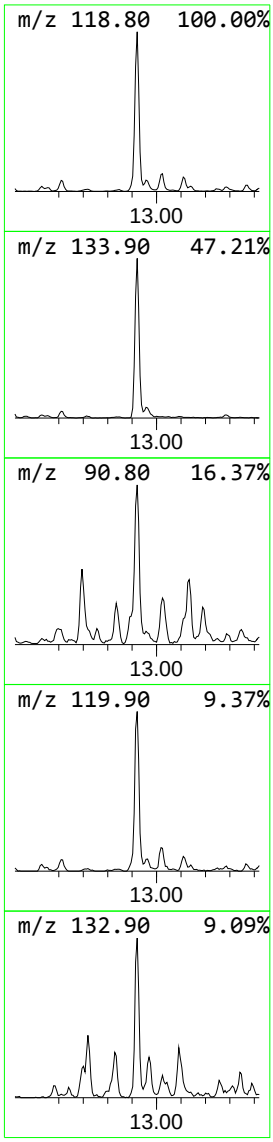
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.920	18.73 ug/l	873905	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	93
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

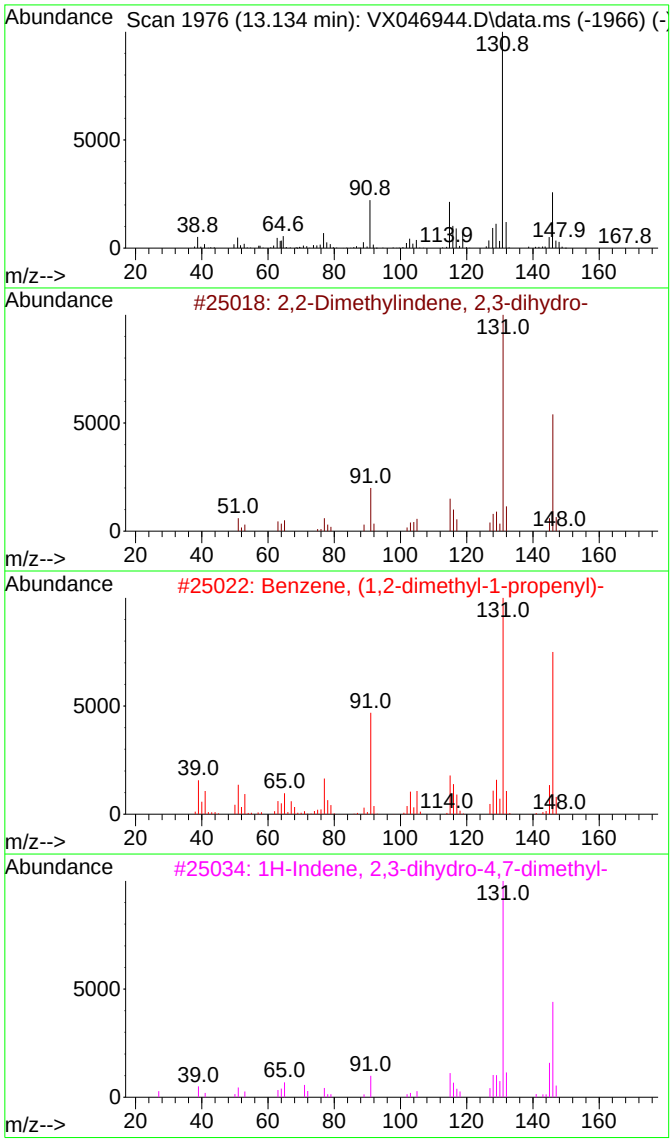
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.M
Quant Title : SW846 8260

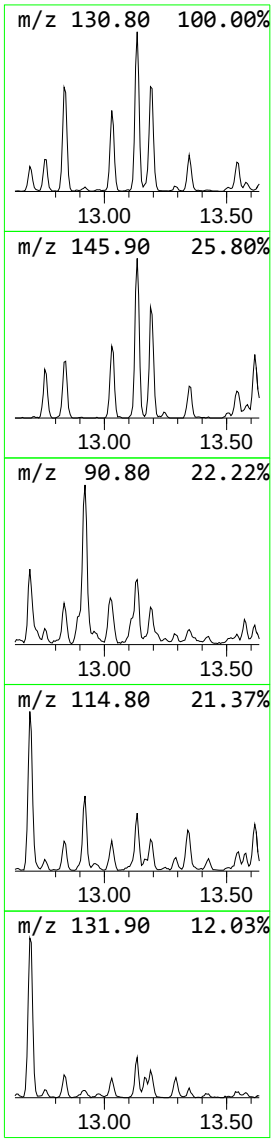
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 2,2-Dimethylindene, 2,3-dih... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.134	7.40 ug/l	345077	1,4-Dichlorobenzene-d4	12.018

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
2	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91
4	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	90
5	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071025\
Data File : VX046944.D
Acq On : 10 Jul 2025 15:04
Operator : JC/MD
Sample : Q2553-02
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
AOC-202

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X070225W.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
Butane, 2-methyl-	1.770	13.1	ug/l	349338	1	5.568	1335390	50.0
Butane, 2,2-dim...	2.361	7.0	ug/l	188371	1	5.568	1335390	50.0
Butane, 2,3-dim...	2.800	23.3	ug/l	622395	1	5.568	1335390	50.0
Pentane, 3-methyl-	3.117	15.7	ug/l	418139	1	5.568	1335390	50.0
Cyclopentane, 1...	5.861	5.8	ug/l	156343	1	5.568	1335390	50.0
Cyclopentene, 1...	9.165	7.8	ug/l	355664	3	10.055	2288240	50.0
Benzene, 1,4-di...	12.244	15.4	ug/l	719008	4	12.018	2332730	50.0
Benzene, 1-ethe...	12.695	10.7	ug/l	499998	4	12.018	2332730	50.0
Benzene, 1,2,3,...	12.920	18.7	ug/l	873905	4	12.018	2332730	50.0
2,2-Dimethylind...	13.134	7.4	ug/l	345077	4	12.018	2332730	50.0