LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

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_N\methods\82N060625W.M
Title : SW846 8260

Signal : TIC: VN087115.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.354	62	68	73	rBV2	9270	15632	1.40%	0.149%
2	2.554	93	102	109	rBV	19539	35730	3.19%	0.341%
3	3.283	216	226	240	rBV	183371	455563	40.72%	4.344%
4	3.671	286	292	301	rVB2	6269	14911	1.33%	0.142%
5	4.448	412	424	441	rVB3	33476	121290	10.84%	1.157%
6	5.306	554	570	581	rBV2	71641	303134	27.10%	2.890%
7	5.836	646	660	677	rBV2	56375	197958	17.70%	1.888%
8	6.818	815	827	837	rVB6	5241	14193	1.27%	0.135%
9	7.089	863	873	881	rBV4	6110	17985	1.61%	0.171%
10	7.236	889	898	907	rBV2	22779	68310	6.11%	0.651%
11	7.341	907	916	923	rBV2	29768	82378	7.36%	0.785%
12	8.012	1024	1030	1040	rVB4	9469	24256	2.17%	0.231%
13	8.171	1048	1057	1062	rBV	140459	338778	30.28%	3.230%
14	8.230	1062	1067	1078	rVV	225382	548512	49.03%	5.230%
15	8.335	1078	1085	1096	rVB2	62468	155050	13.86%	1.478%
16	8.518	1109	1116	1120	rBV4	8880	21837	1.95%	0.208%
17	8.588	1120	1128	1139	rVV	153634	359488	32.14%	3.428%
18	8.724	1139	1151	1153	rVV5	12266	33624	3.01%	0.321%
19	8.765	1153	1158	1169	rVV3	30606	100525	8.99%	0.959%
20	8.859	1169	1174	1182	rVB5	10965	24496	2.19%	0.234%
21	9.106	1207	1216	1228	rBV	381737	805248	71.98%	7.678%
22	9.335	1250	1255	1260	rBV3	6251	12087	1.08%	0.115%
23	9.565	1287	1294	1304	rBV5	20975	60878	5.44%	0.580%
24	9.865	1340	1345	1351	rVB2	8114	15972	1.43%	0.152%
25	10.018	1364	1371	1378	rVB2	23139	49433	4.42%	0.471%
26	10.112	1380	1387	1388	rBV3	8885	14970	1.34%	0.143%
27	10.147	1388	1393	1406	rVB	45031	95277	8.52%	0.908%
28	10.329	1418	1424	1431	rVB5	9432	18377	1.64%	0.175%
29	10.453	1439	1445	1448	rBV4	8714	14753	1.32%	0.141%
30	10.494	1448	1452	1457	rVB3	8184	14444	1.29%	0.138%
31	10.565	1457	1464	1480	rBV	587987	1118665	100.00%	10.667%
32	10.759	1485	1497	1504	rVV4	33169	73967	6.61%	0.705%
33	10.829	1504	1509	1517	rVV3	28059	54114	4.84%	0.516%
34	10.912	1517	1523	1529	rVV3	9273	17137	1.53%	0.163%
35	10.982	1529	1535	1545	rVV5	23554	64406	5.76%	0.614%
36	11.253	1574	1581	1586	rBV4	11728	23630	2.11%	0.225%

LSC Area Percent Report

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Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

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_N\methods\82N060625W.M
Title : SW846 8260

37	11.382	1599	1603	1609	rVB4	6852	12062	1.08%	0.115%
38	11.865	1677	1685	1694	rBV	498496	914681	81.77%	8.722%
39	11.965	1698	1702	1707	rVB5	7585	12855	1.15%	0.123%
40	12.076	1713	1721	1724	rBV3	10920	20691	1.85%	0.197%
41	12.582	1792	1807	1810	rBV7	4584	13167	1.18%	0.126%
42	12.694	1821	1826	1833	rVB	46256	76019	6.80%	0.725%
43	12.847	1845	1852	1862	rBV	382463	661835	59.16%	6.311%
44	13.029	1874	1883	1889	rBV	12430	24637	2.20%	0.235%
45	13.335	1929	1935	1942	rBV2	9515	18157	1.62%	0.173%
46	13.482	1955	1960	1967	rVB	26552	44173	3.95%	0.421%
47	13.600	1973	1980	1989	rBV2	29746	69409	6.20%	0.662%
48	13.788	2005	2012	2024	rBV	514553	917060	81.98%	8.744%
49	13.882	2024	2028	2033	rVB2	18413	31331	2.80%	0.299%
50	13.947	2033	2039	2043	rBV	32701	52609	4.70%	0.502%
51	13.994	2043	2047	2049	rVV	31816	46279	4.14%	0.441%
52	14.029	2049	2053	2062	rVB	79228	131643	11.77%	1.255%
53	14.112	2062	2067	2073	rVB	47016	73898	6.61%	0.705%
54	14.176	2073	2078	2084	rVB4	10004	17416	1.56%	0.166%
55	14.270	2085	2094	2098	rBV2	32107	58682	5.25%	0.560%
56	14.329	2098	2104	2109	rVB	47432	75483	6.75%	0.720%
57	14.441	2113	2123	2130	rBV	394852	676462	60.47%	6.450%
58	14.512	2130	2135	2138	rBV4	10936	21774	1.95%	0.208%
59	14.576	2143	2146	2151	rVV2	21999	36265	3.24%	0.346%
60	14.653	2151	2159	2169	rVB	102149	196406	17.56%	1.873%
61	14.770	2174	2179	2182	rBV	23702	37257	3.33%	0.355%
62	14.876	2190	2197	2201	rVB3	13512	30404	2.72%	0.290%
63	14.935	2201	2207	2212	rBV4	33192	64401	5.76%	0.614%
64	15.117	2232	2238	2243	rBV2	42169	80586	7.20%	0.768%
65	15.200	2248	2252	2257	rVV3	10919	18619	1.66%	0.178%
66	15.253	2257	2261	2265	rVB5	8867	12977	1.16%	0.124%
67	15.317	2265	2272	2277	rBV3	33728	66004	5.90%	0.629%
68	15.382	2277	2283	2284	rVV2	22432	38250	3.42%	0.365%
69	15.423	2285	2290	2298	rVB4	71622	144308	12.90%	1.376%
70	15.594	2313	2319	2324	rVB3	20281	36859	3.29%	0.351%
71	15.682	2330	2334	2341	rBV7	11258	23592	2.11%	0.225%
72	15.794	2347	2353	2360	rVB4	17826	48836	4.37%	0.466%
73	16.059	2388	2398	2407	rBV6	16239	60005	5.36%	0.572%
74	16.153	2409	2414	2421	rVV5	13851	37487	3.35%	0.357%
75	16.253	2425	2431	2438	rVV	30745	71410	6.38%	0.681%
76	16.335	2438	2445	2455	rVB6	20754	57185	5.11%	0.545%

LSC Area Percent Report

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Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

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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Peak separation: 5

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Title : SW846 8260

77	16.482	2464	2470	2476	rBV8	8376	21489	1.92%	0.205%
78	16.670	2496	2502	2503	rBV3	10183	14685	1.31%	0.140%
79	16.711	2505	2509	2515	rVB2	18644	33143	2.96%	0.316%

Sum of corrected areas: 10487499

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Data File : VN087115.D
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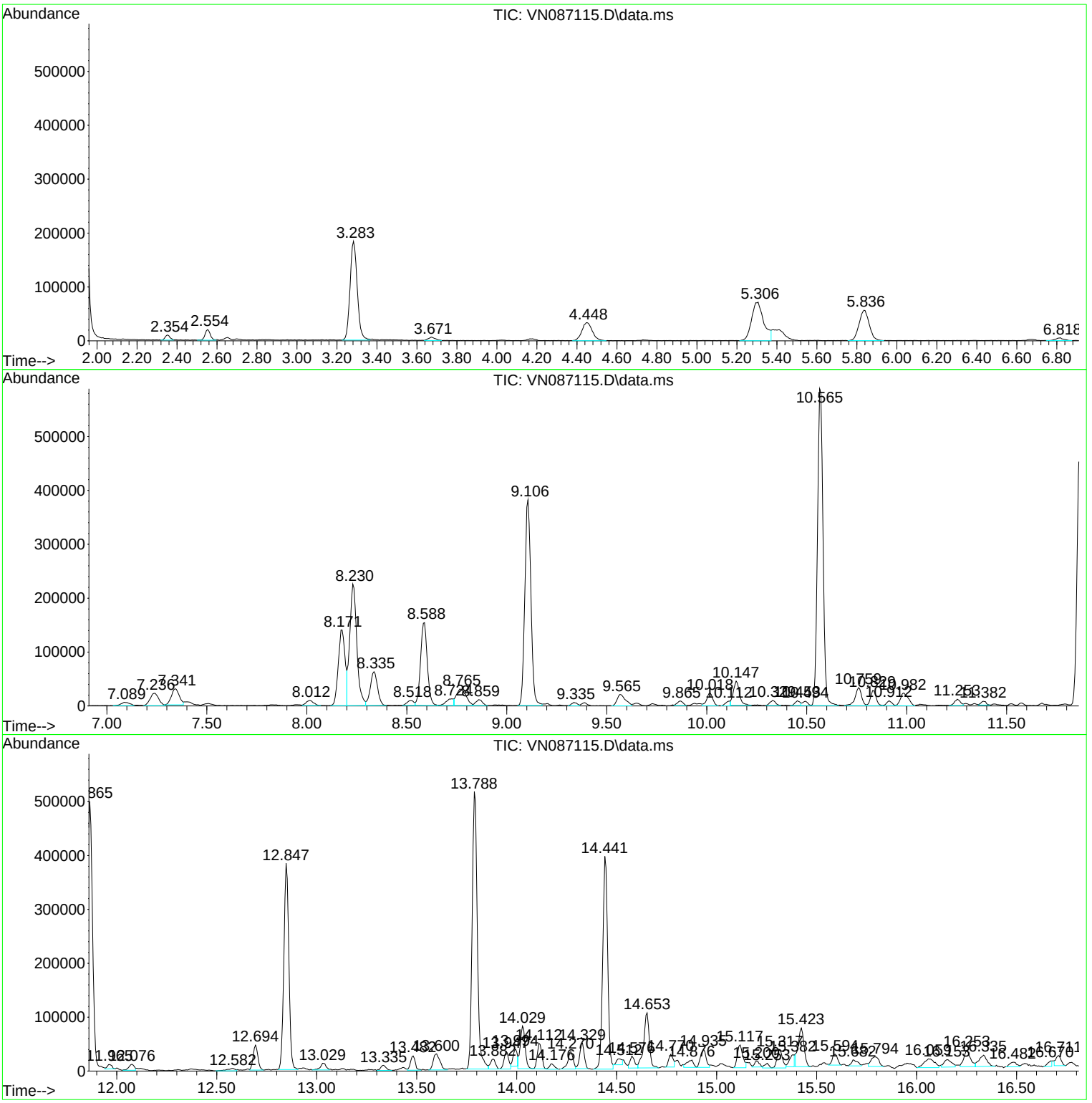
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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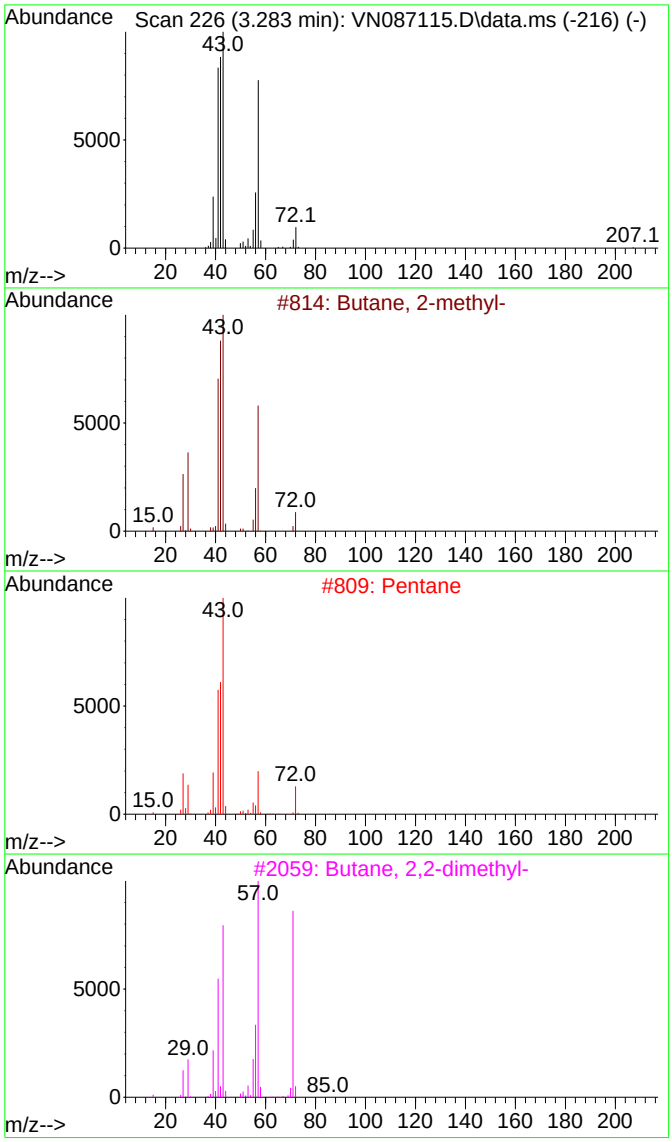
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

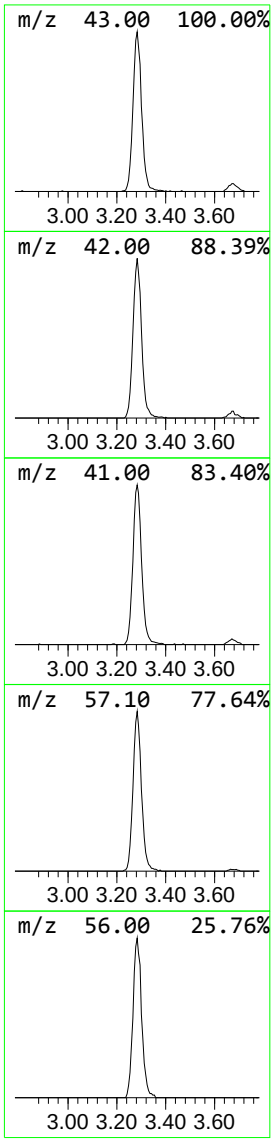
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.283	41.53 ug/l	455563	Pentafluorobenzene	8.230

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	38
3	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	38
4	2-Pentanone, 3-methyl-	100	C6H12O	000565-61-7	17
5	1,1-Diisobutoxy-butane	202	C12H26O2	013002-16-9	16





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

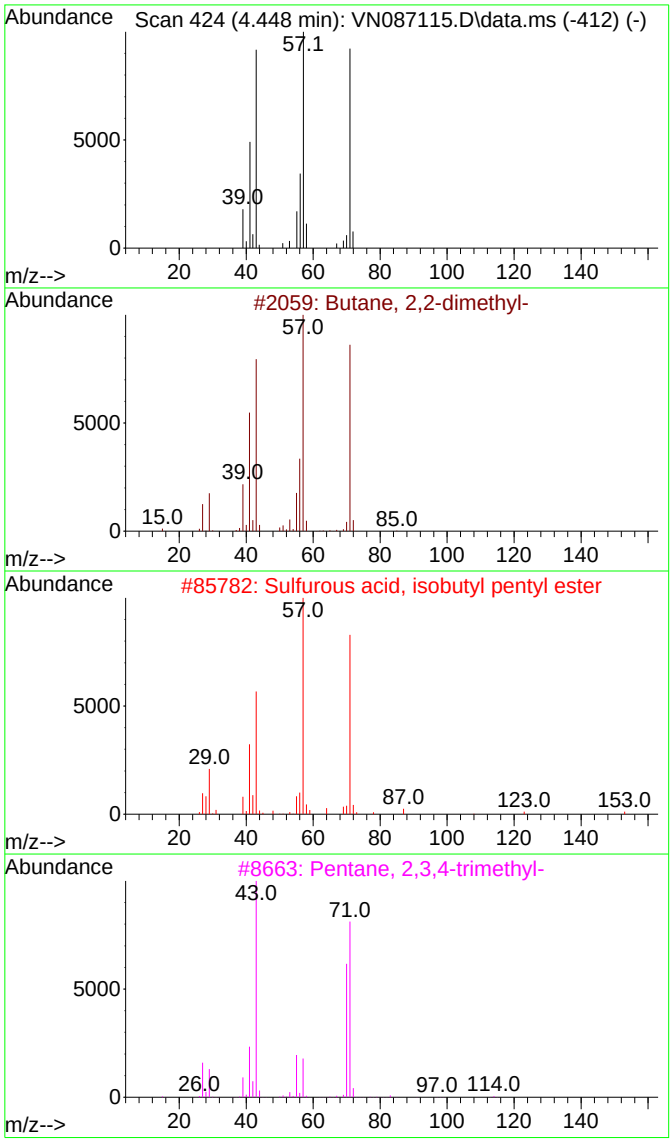
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

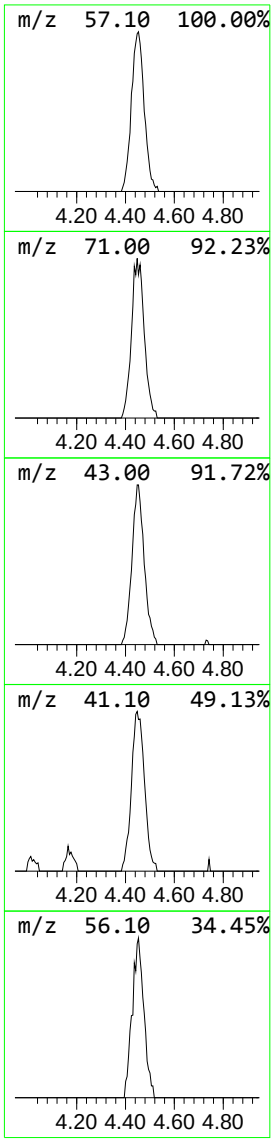
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.448	11.06 ug/l	121290	Pentafluorobenzene	8.230

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	50
3	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	36
4	Pentane, 3-bromo-	150	C5H11Br	001809-10-5	33
5	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	33





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

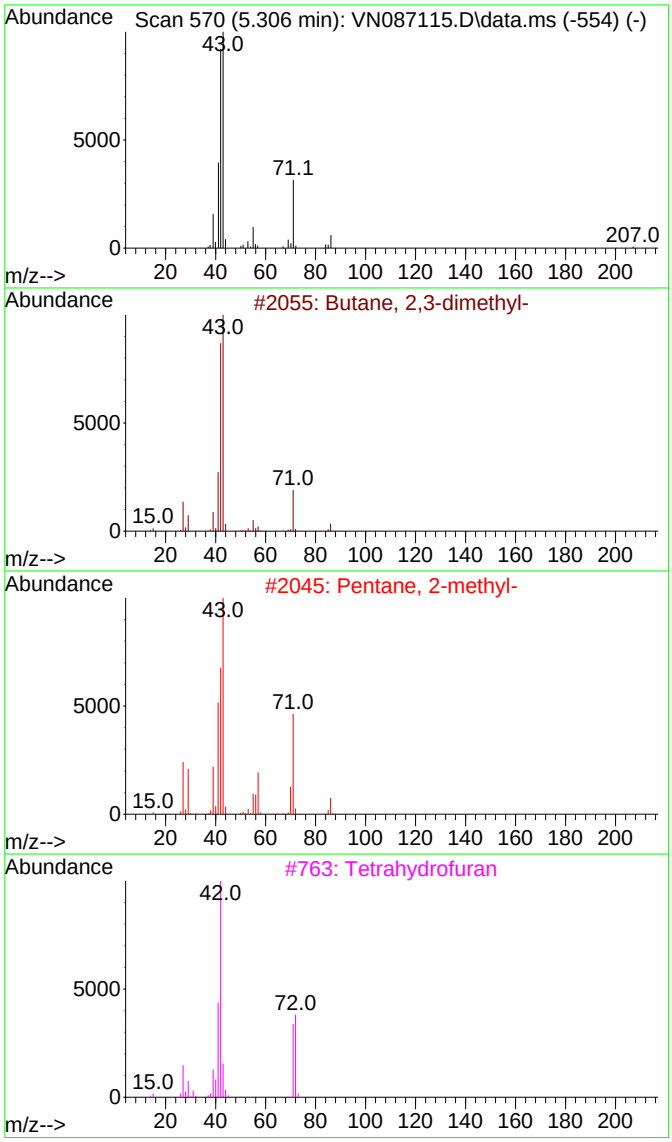
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

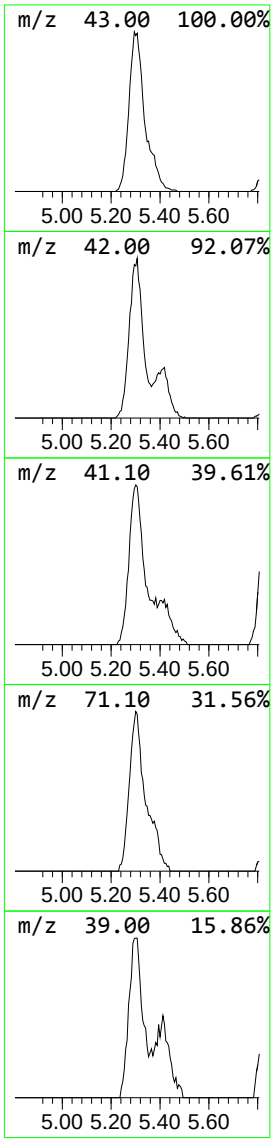
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.306	27.63 ug/l	303134	Pentafluorobenzene	8.230

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	52
3	Tetrahydrofuran	72	C4H8O	000109-99-9	38
4	Pentane	72	C5H12	000109-66-0	38
5	1,2,4,5-Tetrazine, 1,4-dihydro-3...	112	C4H8N4	037454-64-1	28





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Instrument :
MSVOA_N
ClientSampleId :
MW-10

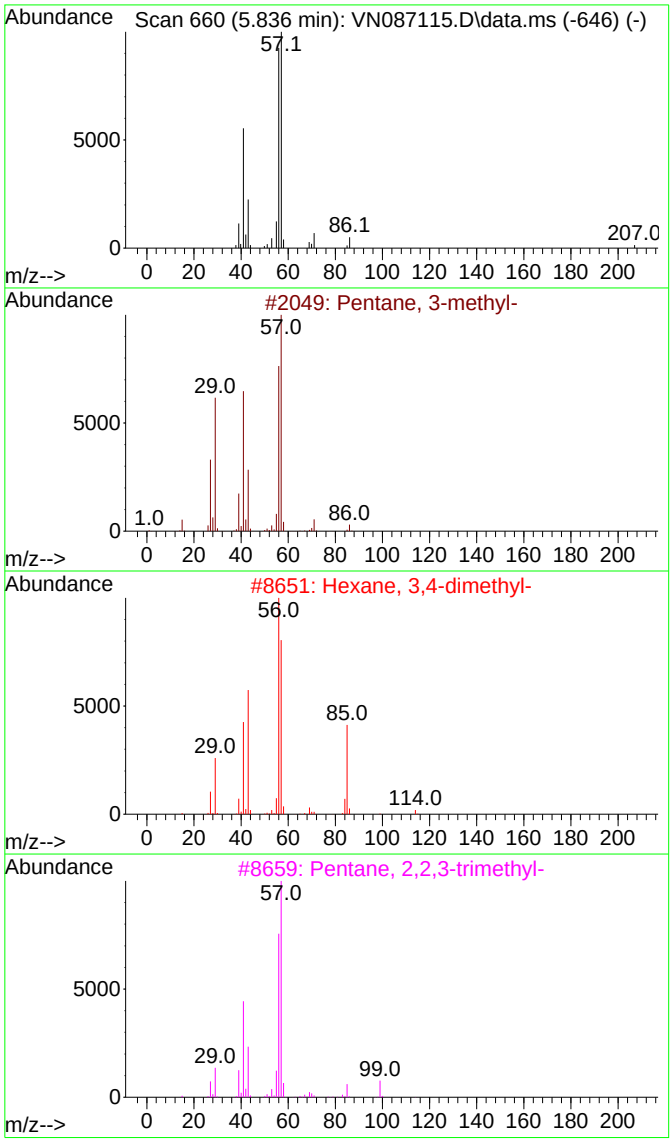
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

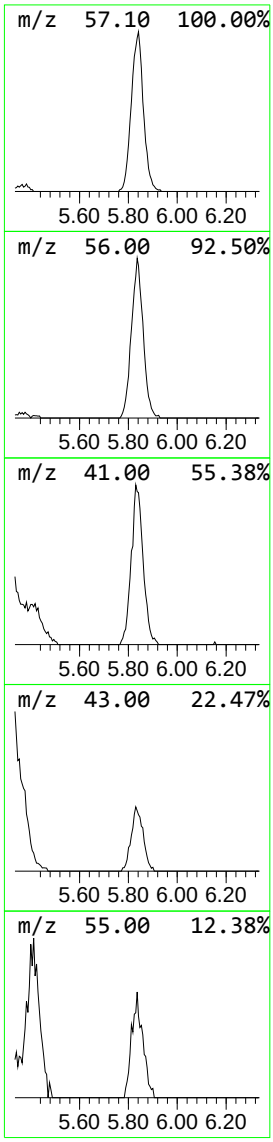
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.836	18.05 ug/l	197958	Pentafluorobenzene	8.230

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 3,4-dimethyl-	114	C8H18	000583-48-2	78
3	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	72
4	Sulphuric acid dibutyl ester	210	C8H18O4S	000625-22-9	56
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	56





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Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

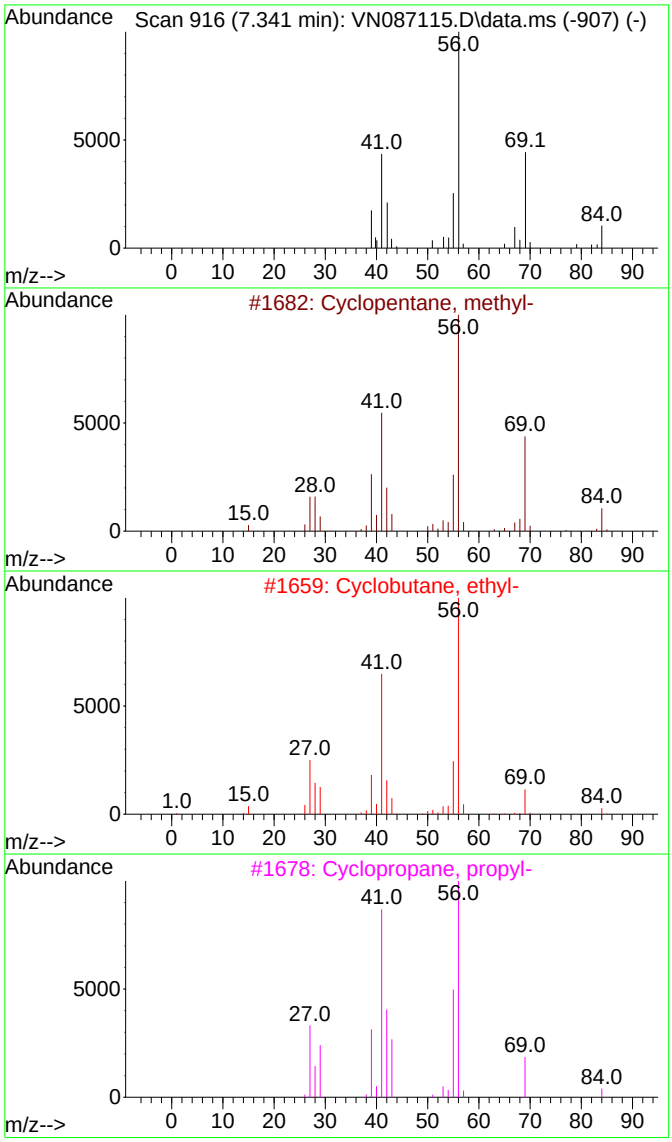
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

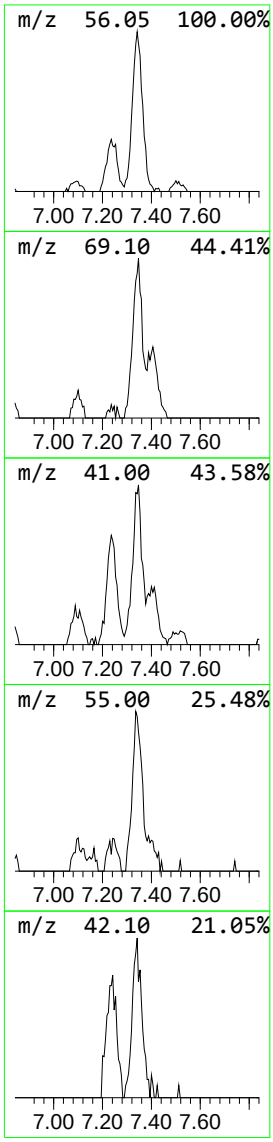
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.341	7.51 ug/l	82378	Pentafluorobenzene	8.230

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
3	Cyclopropane, propyl-	84	C6H12	002415-72-7	64
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5	1-Pentanol, 3-methyl-	102	C6H14O	000589-35-5	38





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Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

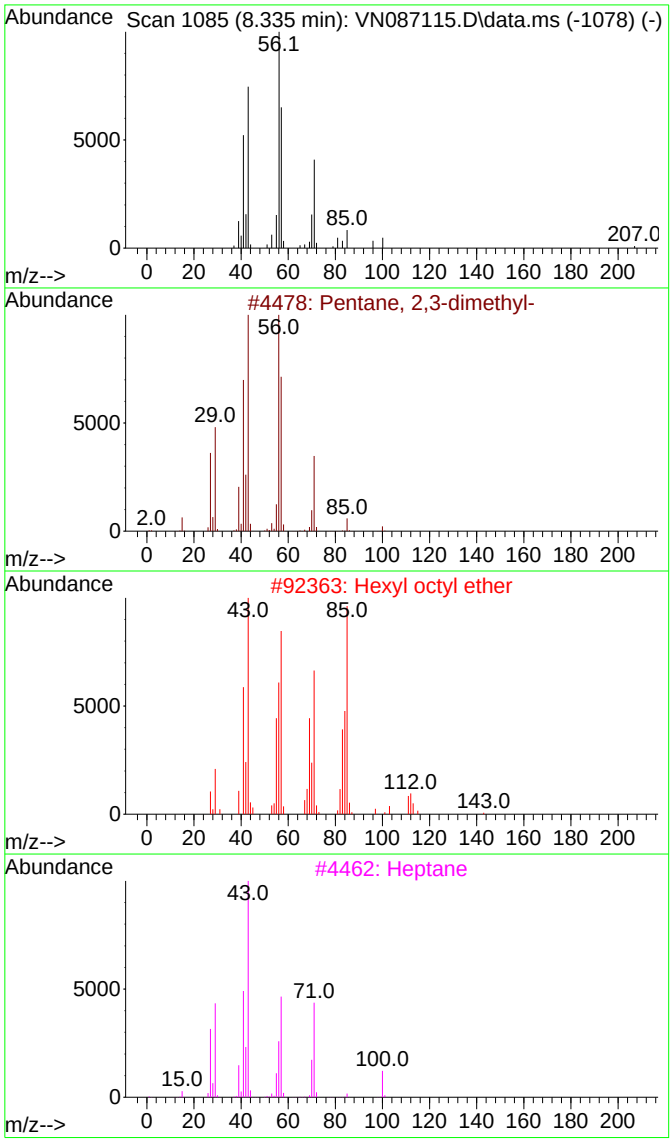
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

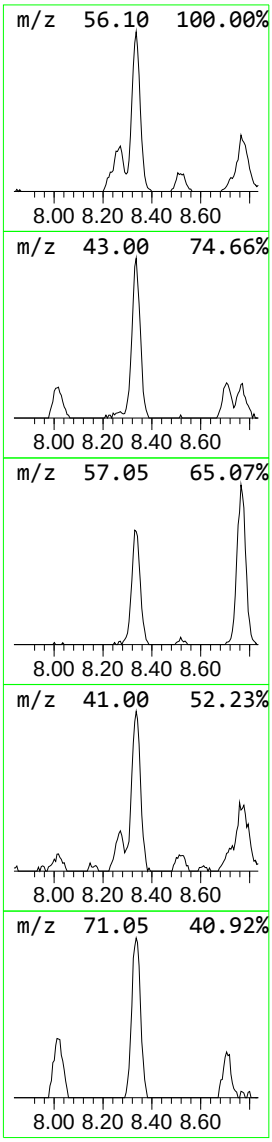
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.335	14.13 ug/l	155050	Pentafluorobenzene	8.230

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	78
2	Hexyl octyl ether	214	C14H30O	017071-54-4	43
3	Heptane	100	C7H16	000142-82-5	43
4	1-Hexanol, 5-methyl-2-(1-methyle...	158	C10H22O	002051-33-4	43
5	Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

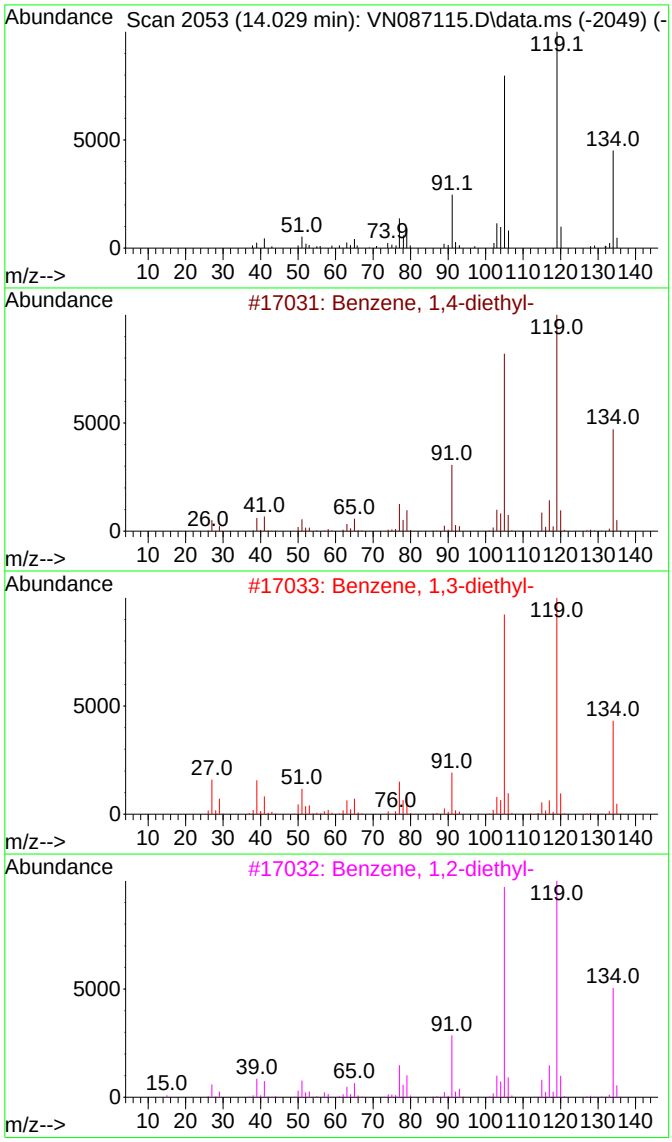
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

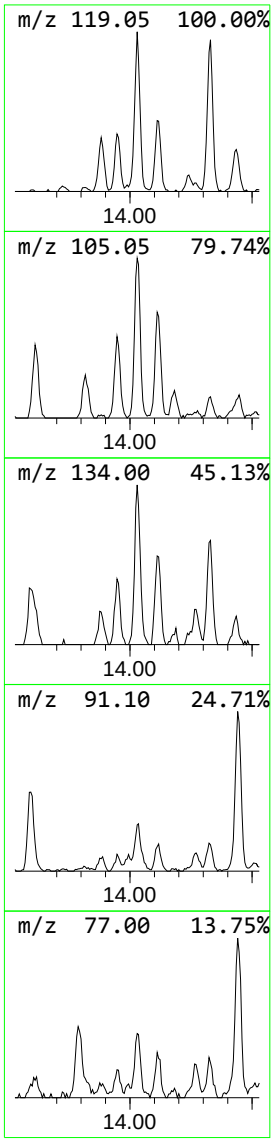
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.029	7.18 ug/l	131643	1,4-Dichlorobenzene-d4	13.788

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
2	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
3	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	91
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	80
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

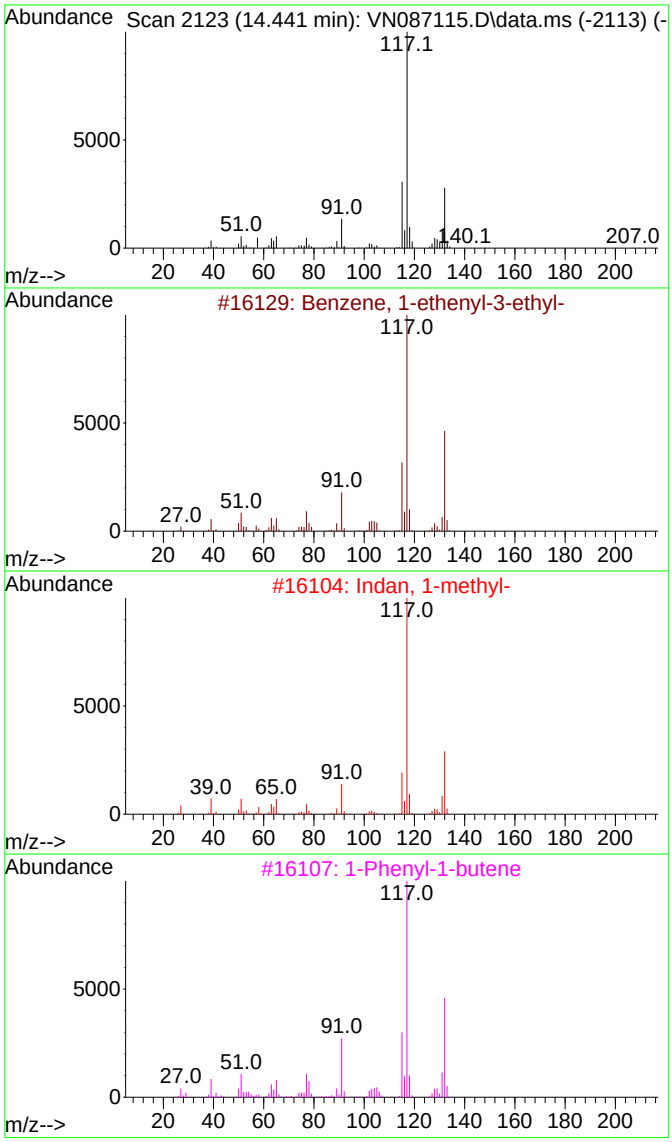
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

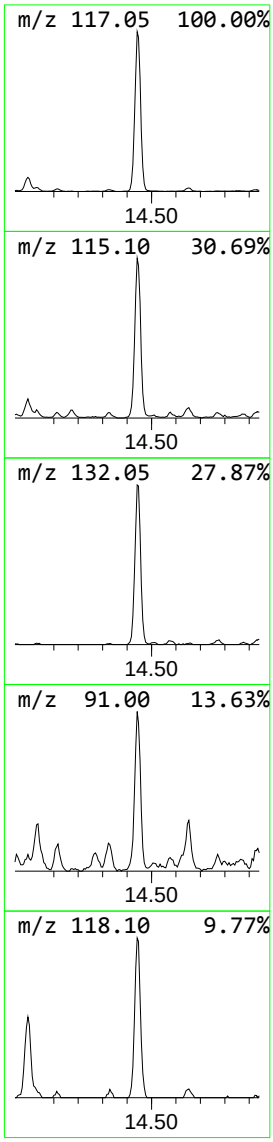
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.441	36.88 ug/l	676462	1,4-Dichlorobenzene-d4	13.788

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	87
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

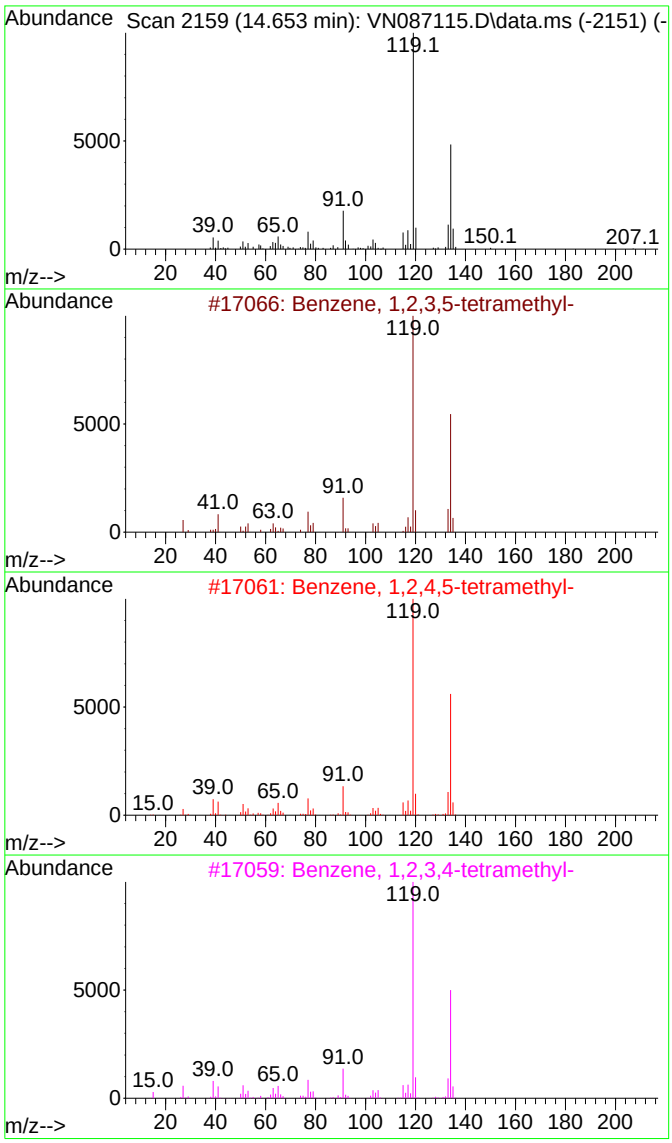
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

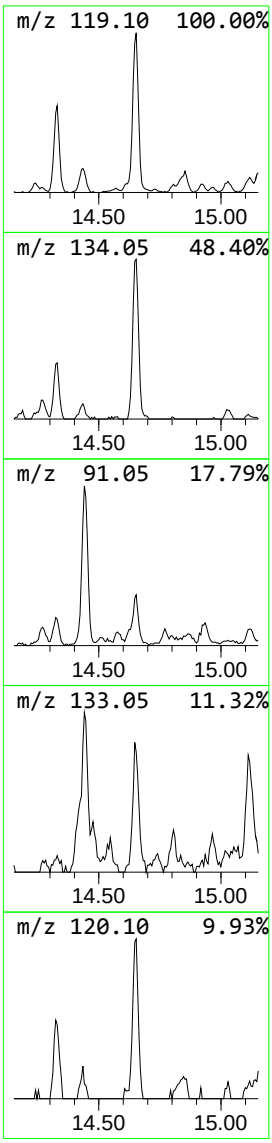
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.653	10.71 ug/l	196406	1,4-Dichlorobenzene-d4	13.788

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5	1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	91





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Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
Sample : Q2333-03
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

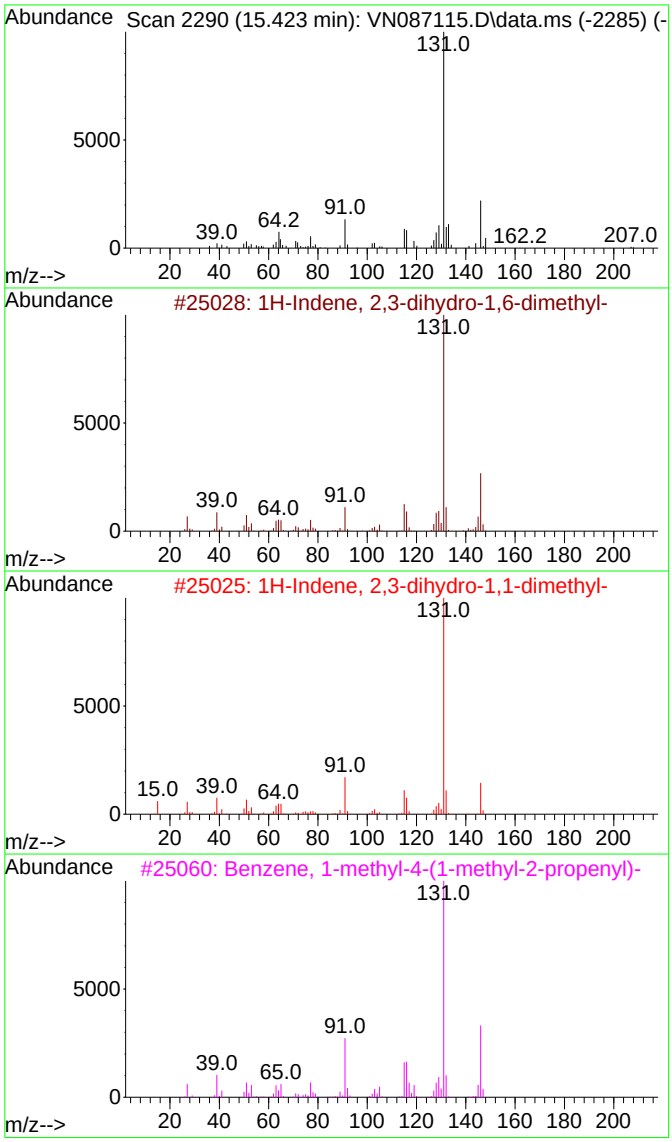
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.M
Quant Title : SW846 8260

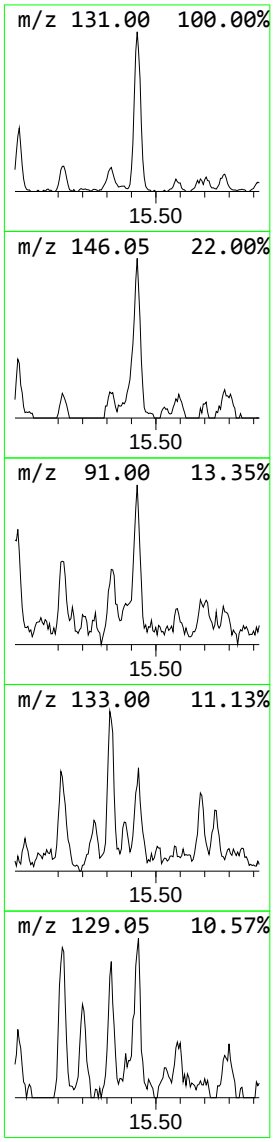
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.423	7.87 ug/l	144308	1,4-Dichlorobenzene-d4	13.788

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	90
2	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	83
3	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	80
4	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	72
5	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061925\
Data File : VN087115.D
Acq On : 19 Jun 2025 16:28
Operator : JC\MD
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Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 17 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N060625W.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-	3.283	41.5	ug/l	455563	1	8.230	548512	50.0
Butane, 2,2-dim...	4.448	11.1	ug/l	121290	1	8.230	548512	50.0
Butane, 2,3-dim...	5.306	27.6	ug/l	303134	1	8.230	548512	50.0
Pentane, 3-methyl-	5.836	18.1	ug/l	197958	1	8.230	548512	50.0
Cyclopentane, m...	7.341	7.5	ug/l	82378	1	8.230	548512	50.0
Pentane, 2,3-di...	8.335	14.1	ug/l	155050	1	8.230	548512	50.0
Benzene, 1,4-di...	14.029	7.2	ug/l	131643	4	13.788	917060	50.0
Benzene, 1-ethe...	14.441	36.9	ug/l	676462	4	13.788	917060	50.0
Benzene, 1,2,3,...	14.653	10.7	ug/l	196406	4	13.788	917060	50.0
1H-Indene, 2,3-...	15.423	7.9	ug/l	144308	4	13.788	917060	50.0