LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

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\82N021825W.M
Title : SW846 8260

Signal : TIC: VN085924.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.236	38	48	65	rVB	241696	559127	46.04%	2.422%
2	2.636	107	116	128	rVB2	65759	157295	12.95%	0.681%
3	3.418	236	249	261	rBV3	33378	113936	9.38%	0.493%
4	4.283	377	396	400	rBV3	103491	395173	32.54%	1.712%
5	4.348	401	407	425	rVB3	160671	593681	48.88%	2.571%
6	4.895	480	500	520	rBV3	152427	634459	52.24%	2.748%
7	5.271	554	564	568	rBV4	5269	14390	1.18%	0.062%
8	5.512	592	605	618	rVB3	28869	105209	8.66%	0.456%
9	5.936	666	677	689	rVB4	22386	74443	6.13%	0.322%
10	6.100	694	705	718	rBV3	14016	48042	3.96%	0.208%
11	6.447	750	764	775	rBV5	19637	70339	5.79%	0.305%
12	6.624	782	794	803	rBV3	33638	107868	8.88%	0.467%
13	6.742	803	814	826	rVV3	167781	558075	45.95%	2.417%
14	7.524	936	947	952	rBV6	12703	39956	3.29%	0.173%
15	7.583	952	957	969	rVB4	18060	51053	4.20%	0.221%
16	7.783	980	991	998	rVV2	141350	391289	32.22%	1.695%
17	7.865	998	1005	1015	rVV2	328411	898542	73.98%	3.892%
18	7.971	1015	1023	1035	rVB	143352	411289	33.87%	1.781%
19	8.112	1037	1047	1054	rBV	72379	185105	15.24%	0.802%
20	8.171	1054	1057	1064	rVV4	13619	32243	2.65%	0.140%
21	8.253	1064	1071	1085	rVV	150055	404375	33.30%	1.751%
22	8.418	1089	1099	1102	rVV3	70393	177744	14.64%	0.770%
23	8.477	1103	1109	1119	rVV3	142802	522873	43.05%	2.265%
24	8.571	1119	1125	1136	rVV4	61405	183617	15.12%	0.795%
25	8.677	1138	1143	1155	rVB6	17861	56168	4.62%	0.243%
26	8.859	1164	1174	1188	rBV2	506413	1214495	100.00%	5.260%
27	8.965	1188	1192	1198	rVV2	23493	48727	4.01%	0.211%
28	9.041	1198	1205	1211	rVV3	23451	51252	4.22%	0.222%
29	9.118	1211	1218	1222	rVV2	33868	72221	5.95%	0.313%
30	9.171	1222	1227	1240	rVB4	26532	62896	5.18%	0.272%
31	9.359	1250	1259	1261	rBV2	71660	139705	11.50%	0.605%
32	9.400	1262	1266	1272	rVV3	93705	190634	15.70%	0.826%
33	9.459	1272	1276	1285	rVB4	19257	40266	3.32%	0.174%
34	9.547	1285	1291	1293	rBV2	7647	12715	1.05%	0.055%
35	9.594	1293	1299	1306	rVB3	28241	56064	4.62%	0.243%
36	9.694	1306	1316	1323	rBV7	18004	49924	4.11%	0.216%

LSC Area Percent Report

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Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

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\82N021825W.M
Title : SW846 8260

37	9.771	1323	1329	1330	rBV3	12790	19249	1.58%	0.083%
38	9.853	1337	1343	1353	rVB2	75791	163902	13.50%	0.710%
39	9.953	1353	1360	1363	rBV3	34876	66607	5.48%	0.288%
40	10.000	1363	1368	1376	rVB3	78907	170358	14.03%	0.738%
41	10.118	1383	1388	1394	rVB5	15170	34188	2.81%	0.148%
42	10.200	1394	1402	1410	rBV9	16952	54897	4.52%	0.238%
43	10.324	1417	1423	1428	rBV2	33306	63187	5.20%	0.274%
44	10.371	1428	1431	1436	rVV5	12452	20401	1.68%	0.088%
45	10.447	1436	1444	1457	rVV	626879	1198497	98.68%	5.191%
46	10.559	1458	1463	1466	rVV7	9331	23488	1.93%	0.102%
47	10.647	1470	1478	1485	rVV7	22885	79398	6.54%	0.344%
48	10.712	1485	1489	1492	rVV5	9280	15845	1.30%	0.069%
49	10.812	1501	1506	1513	rBV5	12968	28295	2.33%	0.123%
50	10.894	1513	1520	1529	rVB4	57034	137046	11.28%	0.594%
51	11.171	1560	1567	1572	rBV7	14222	34602	2.85%	0.150%
52	11.312	1586	1591	1595	rVV4	13208	24254	2.00%	0.105%
53	11.512	1621	1625	1631	rVB3	8250	13631	1.12%	0.059%
54	11.618	1635	1643	1650	rBV7	7278	17701	1.46%	0.077%
55	11.812	1667	1676	1686	rBV	548882	1055648	86.92%	4.572%
56	11.912	1686	1693	1704	rVV	430080	743128	61.19%	3.219%
57	12.029	1704	1713	1725	rVB	217078	417473	34.37%	1.808%
58	12.359	1759	1769	1775	rBV2	17123	36612	3.01%	0.159%
59	12.665	1813	1821	1827	rBV	141815	242303	19.95%	1.050%
60	12.824	1840	1848	1858	rVB	415553	736762	60.66%	3.191%
61	13.012	1873	1880	1886	rVV	254183	414116	34.10%	1.794%
62	13.076	1886	1891	1892	rVV	48520	67123	5.53%	0.291%
63	13.106	1892	1896	1901	rVV	151336	265700	21.88%	1.151%
64	13.153	1901	1904	1910	rVB2	50633	77139	6.35%	0.334%
65	13.318	1923	1932	1939	rBV	345614	568728	46.83%	2.463%
66	13.418	1942	1949	1951	rVV2	21142	32706	2.69%	0.142%
67	13.465	1951	1957	1966	rVB	514981	804320	66.23%	3.484%
68	13.594	1971	1979	1985	rVB5	31628	73807	6.08%	0.320%
69	13.659	1985	1990	1996	rVB3	12635	19547	1.61%	0.085%
70	13.776	2003	2010	2022	rBV2	532066	1214125	99.97%	5.259%
71	13.871	2022	2026	2032	rVB	30147	49050	4.04%	0.212%
72	13.941	2032	2038	2041	rBV	176935	299202	24.64%	1.296%
73	13.982	2041	2045	2049	rVV	384024	635226	52.30%	2.751%
74	14.018	2049	2051	2060	rVB2	116774	197519	16.26%	0.856%
75	14.106	2060	2066	2071	rVB2	44344	72228	5.95%	0.313%
76	14.176	2071	2078	2083	rBV3	38922	64544	5.31%	0.280%

LSC Area Percent Report

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Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

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

Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Title : SW846 8260

77	14.235	2083	2088	2091	rBV	51989	92922	7.65%	0.402%
78	14.323	2097	2103	2108	rBV	244655	380515	31.33%	1.648%
79	14.388	2108	2114	2117	rBV	142160	244297	20.12%	1.058%
80	14.435	2117	2122	2130	rVB	513540	837378	68.95%	3.627%
81	14.500	2130	2133	2138	rVB4	10539	15794	1.30%	0.068%
82	14.570	2138	2145	2149	rBV3	38109	63630	5.24%	0.276%
83	14.647	2149	2158	2162	rBV	179014	294185	24.22%	1.274%
84	14.682	2162	2164	2169	rVB	54622	71334	5.87%	0.309%
85	14.723	2169	2171	2175	rVB2	13497	16593	1.37%	0.072%
86	14.870	2192	2196	2199	rBV2	16085	23224	1.91%	0.101%
87	14.918	2199	2204	2211	rBV	209165	351612	28.95%	1.523%
88	15.047	2216	2226	2232	rVV2	477397	959171	78.98%	4.155%
89	15.112	2232	2237	2242	rVV4	32360	69573	5.73%	0.301%
90	15.200	2246	2252	2260	rVV3	33702	68269	5.62%	0.296%
91	15.312	2260	2271	2276	rVV3	75337	185542	15.28%	0.804%
92	15.353	2276	2278	2283	rVV2	23899	39564	3.26%	0.171%
93	15.429	2283	2291	2300	rVB4	55391	143617	11.83%	0.622%
94	15.635	2320	2326	2332	rVB2	43649	82802	6.82%	0.359%
95	15.741	2339	2344	2349	rBV3	16565	33166	2.73%	0.144%
96	15.800	2349	2354	2365	rVB3	15687	34544	2.84%	0.150%
97	15.935	2368	2377	2385	rBV3	26225	53408	4.40%	0.231%
98	16.253	2425	2431	2437	rBV3	9329	19303	1.59%	0.084%
99	16.341	2437	2446	2452	rVB6	10983	29173	2.40%	0.126%

Sum of corrected areas: 23087358

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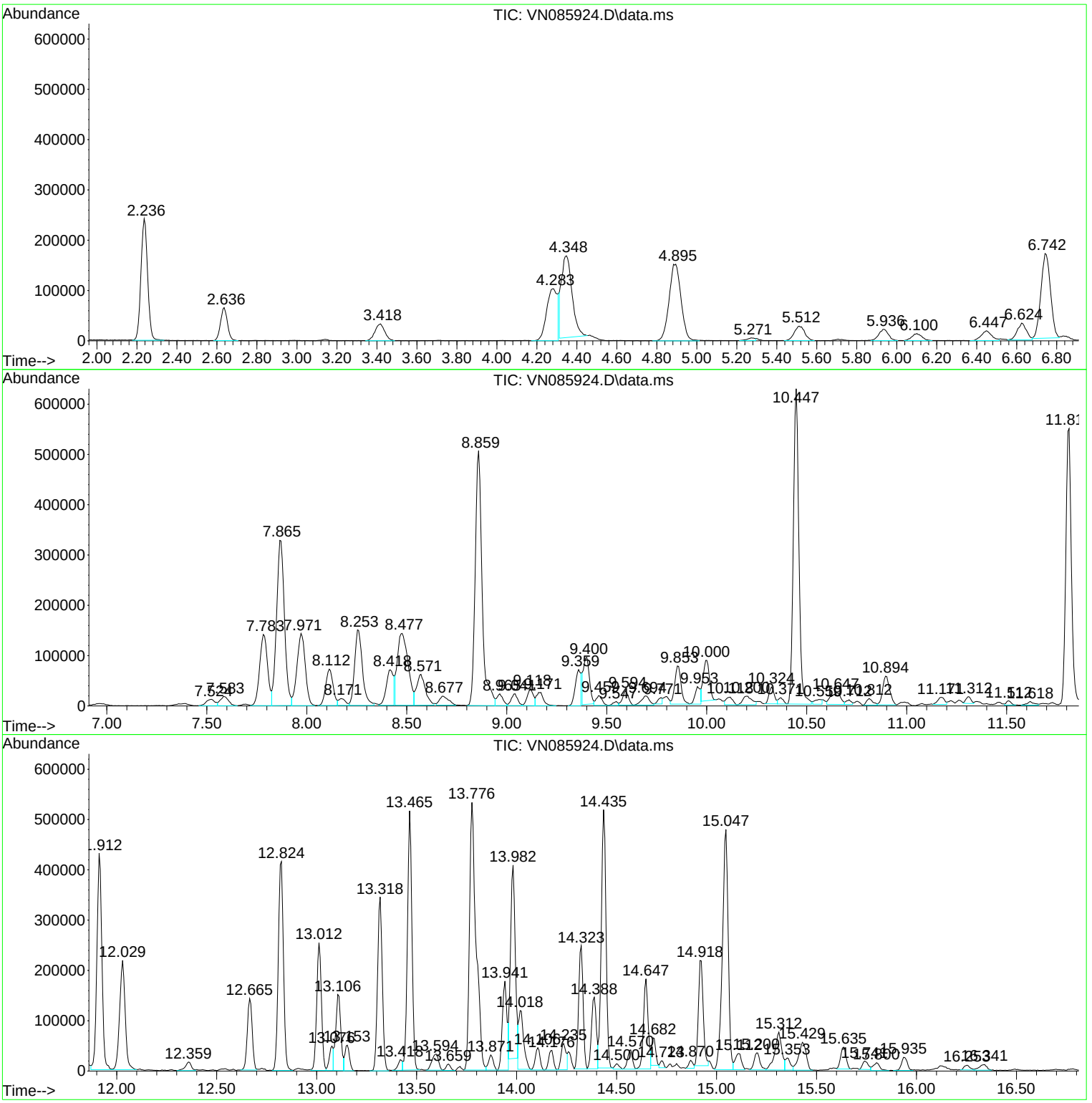
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
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Instrument :
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ClientSampleId :
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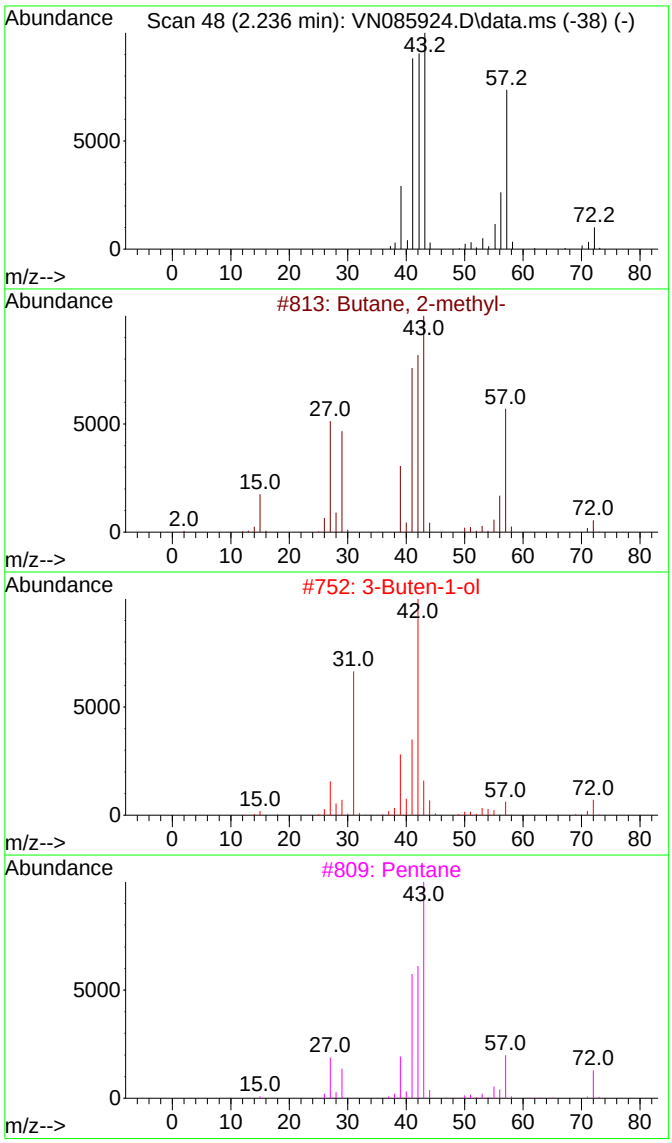
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

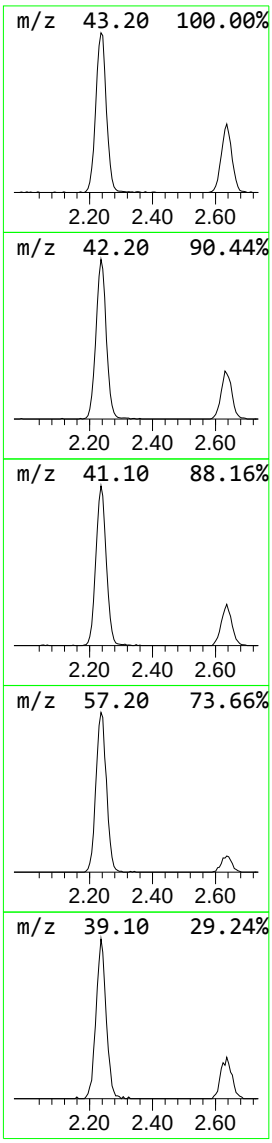
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.236	31.11 ug/l	559127	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	80
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	33
4	1-Pentene	70	C5H10	000109-67-1	25
5	1-Butene	56	C4H8	000106-98-9	10





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

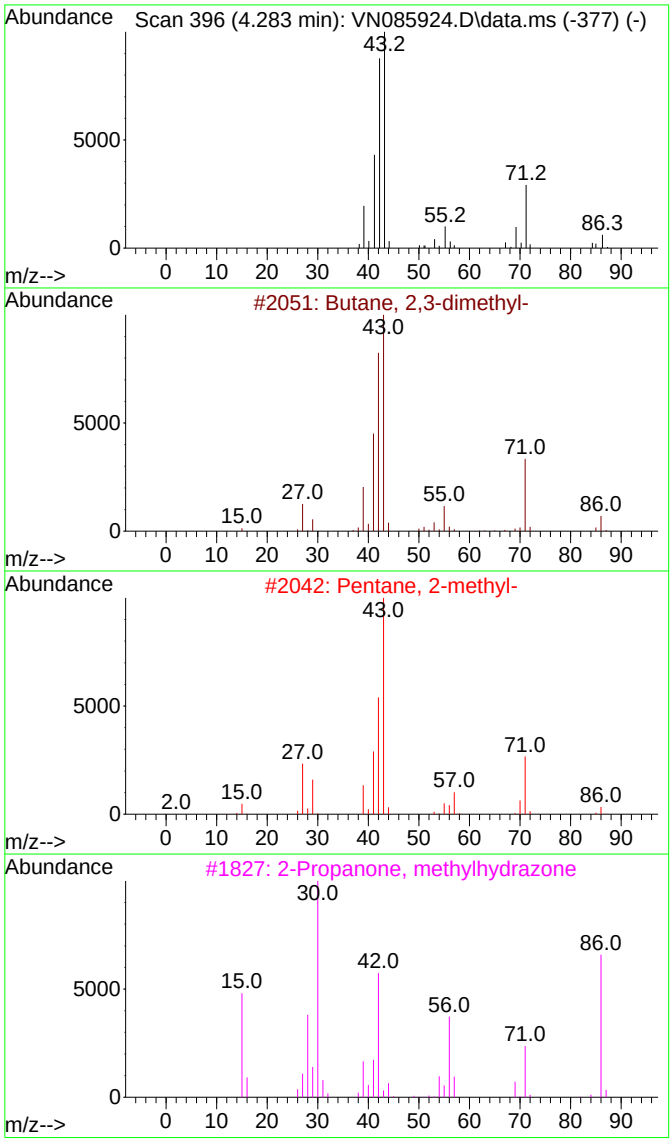
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

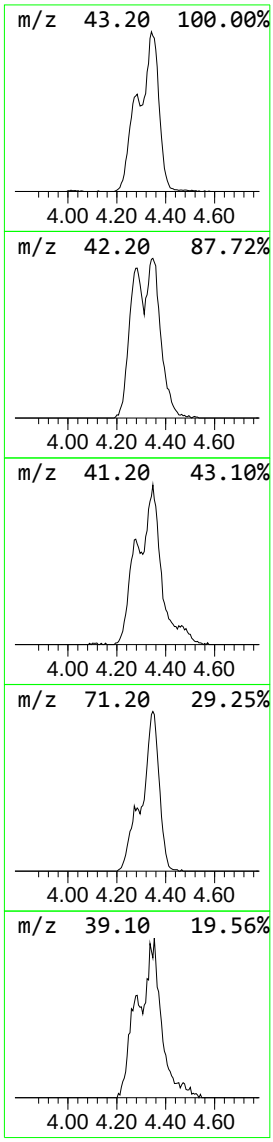
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Butane, 2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.283	21.99 ug/l	395173	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	90
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	47
3	2-Propanone, methylhydrazone	86	C4H10N2	005771-02-8	36
4	1-Pentene	70	C5H10	000109-67-1	33
5	1,2,4,5-Tetrazine, 1,4-dihydro-3...	112	C4H8N4	037454-64-1	28





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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

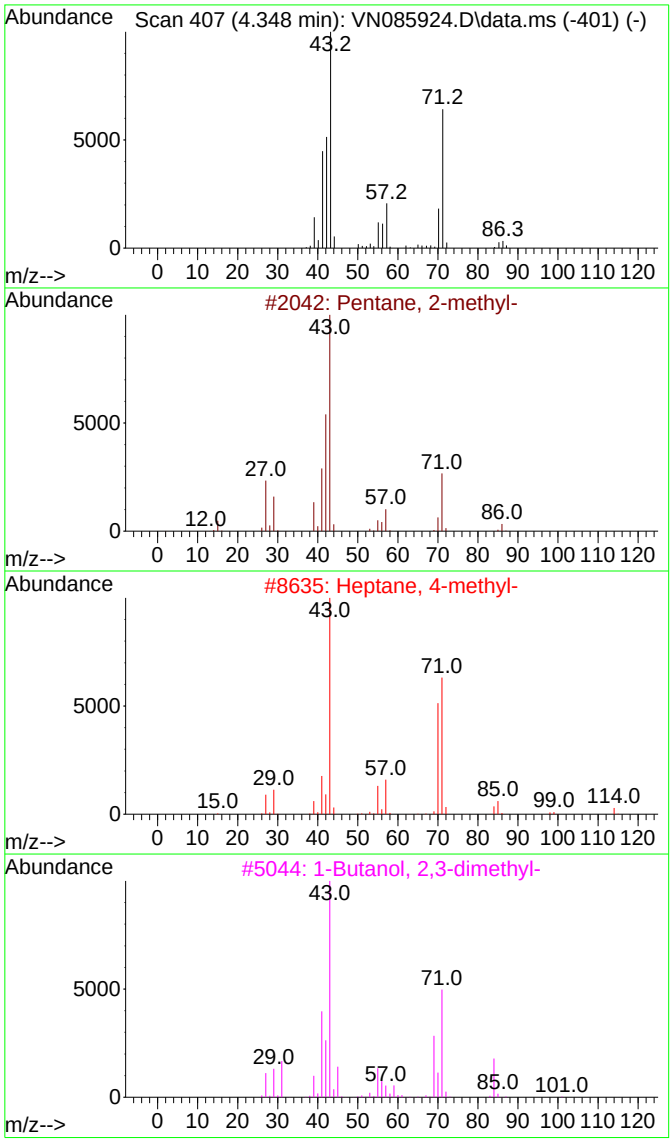
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

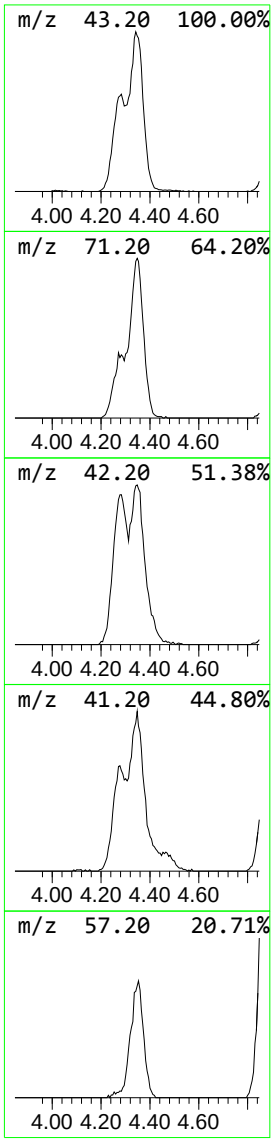
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.348	33.04 ug/l	593681	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	78
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	43
3	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	39
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	37
5	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	33





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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

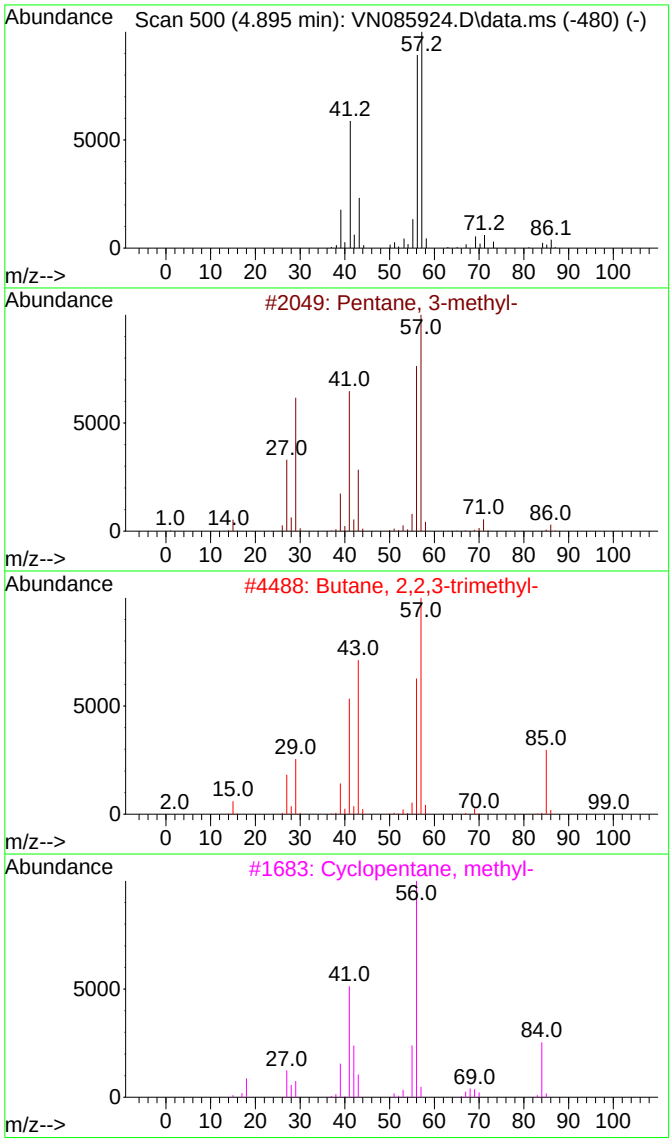
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

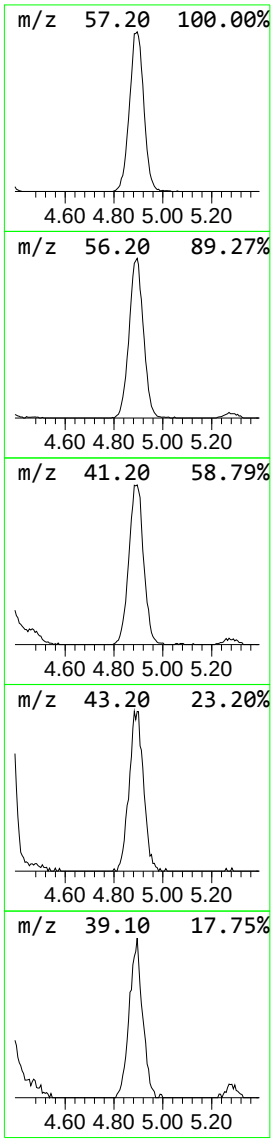
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.895	35.30 ug/l	634459	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	86
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	78
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	58
4	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

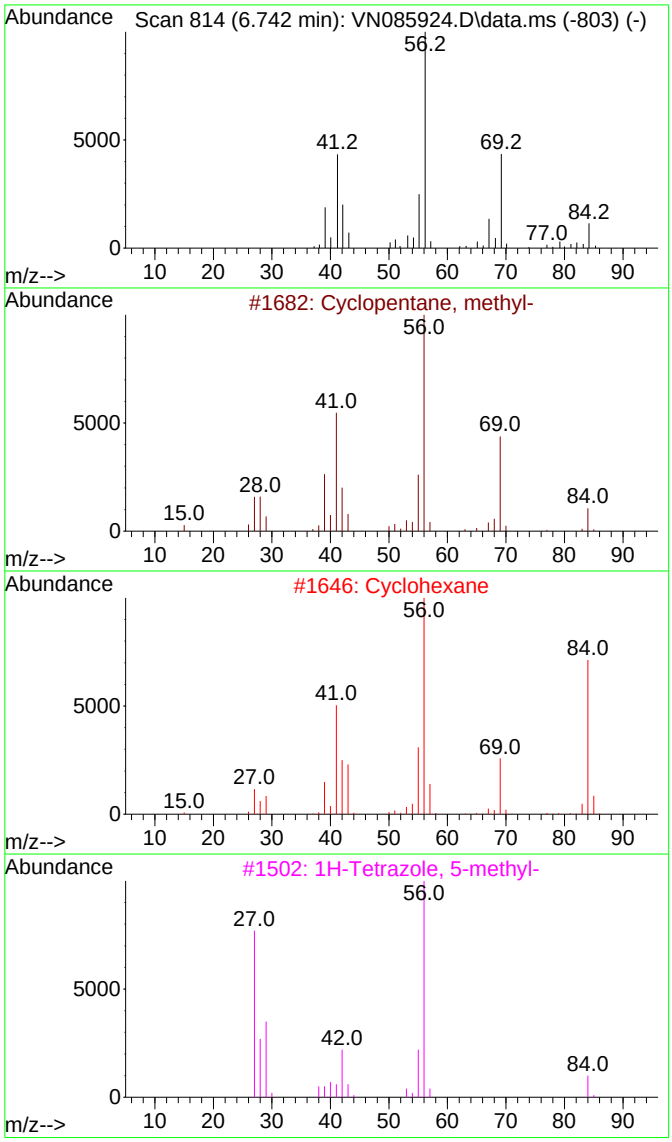
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

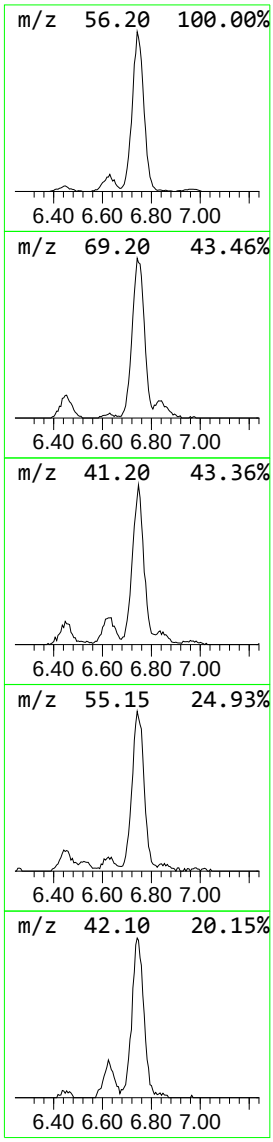
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.742	31.05 ug/l	558075	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclohexane	84	C6H12	000110-82-7	78
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclobutane, ethyl-	84	C6H12	004806-61-5	58
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	43





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Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

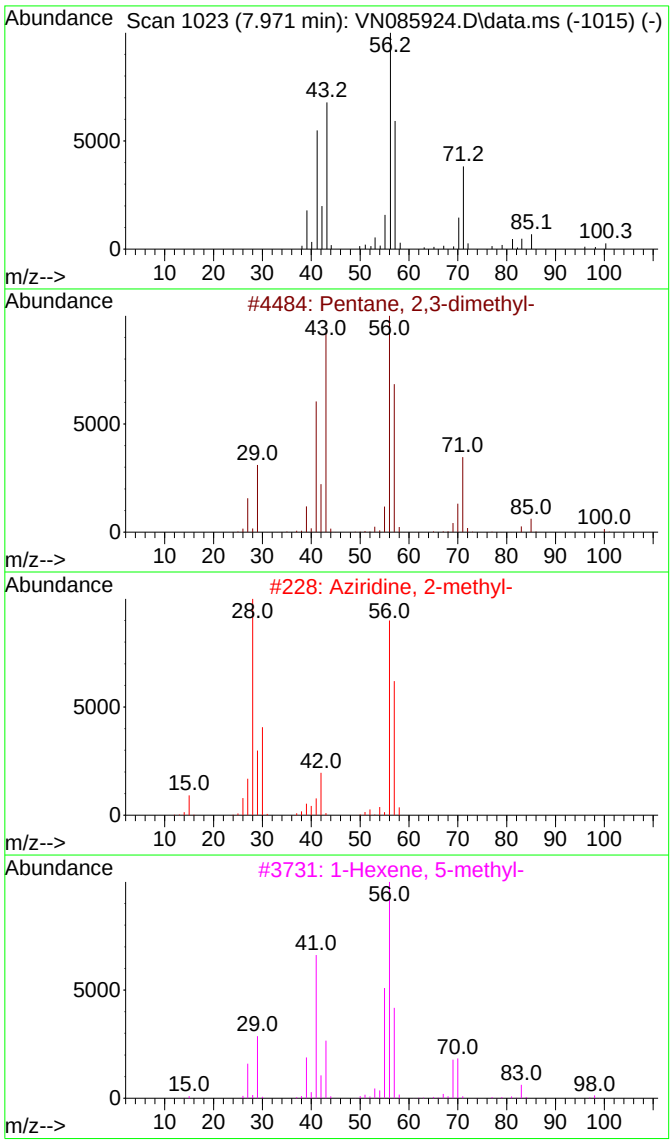
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

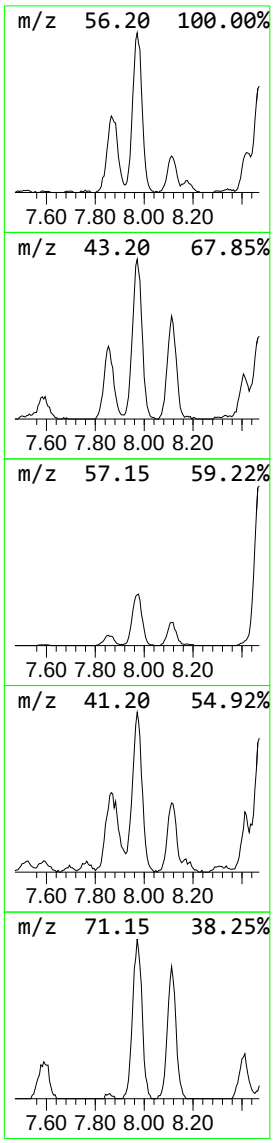
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.971	22.89 ug/l	411289	Pentafluorobenzene	7.871

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	87
2	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	47
3	1-Hexene, 5-methyl-	98	C7H14	003524-73-0	43
4	1,4-Dioxaspiro[2.4]heptan-5-one,...	142	C7H10O3	126091-79-0	42
5	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	38





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Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

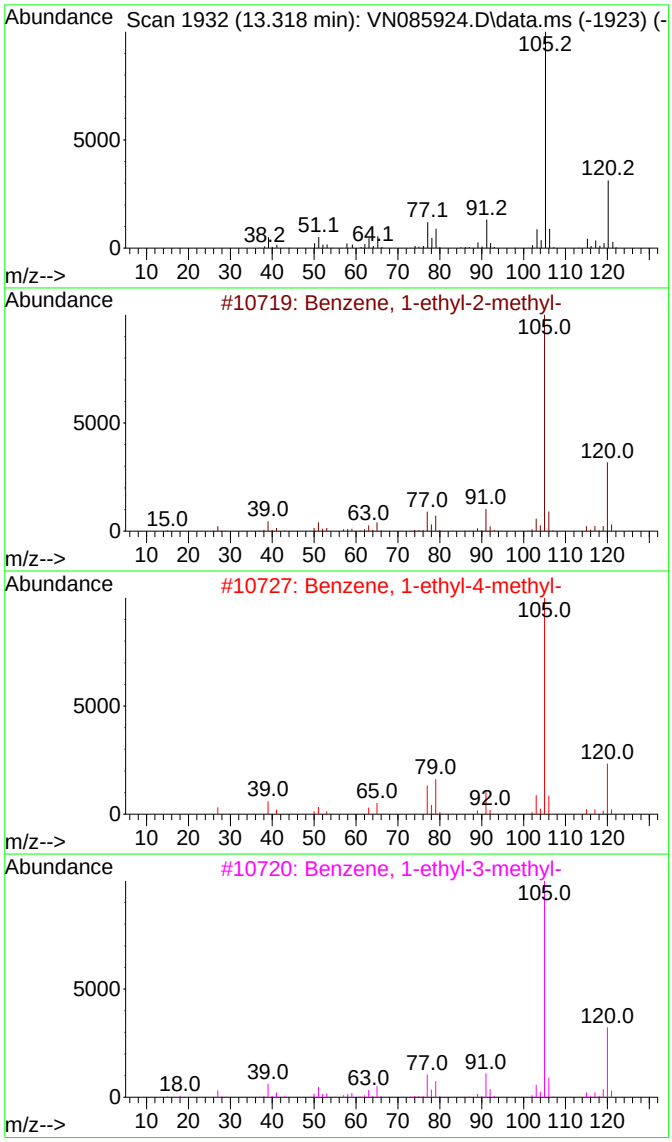
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

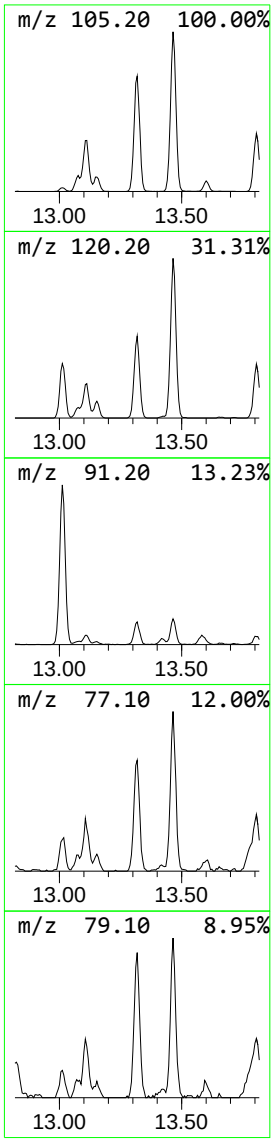
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.318	23.42 ug/l	568728	1,4-Dichlorobenzene-d4	13.776

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	94
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Mesitylene	120	C9H12	000108-67-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





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Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

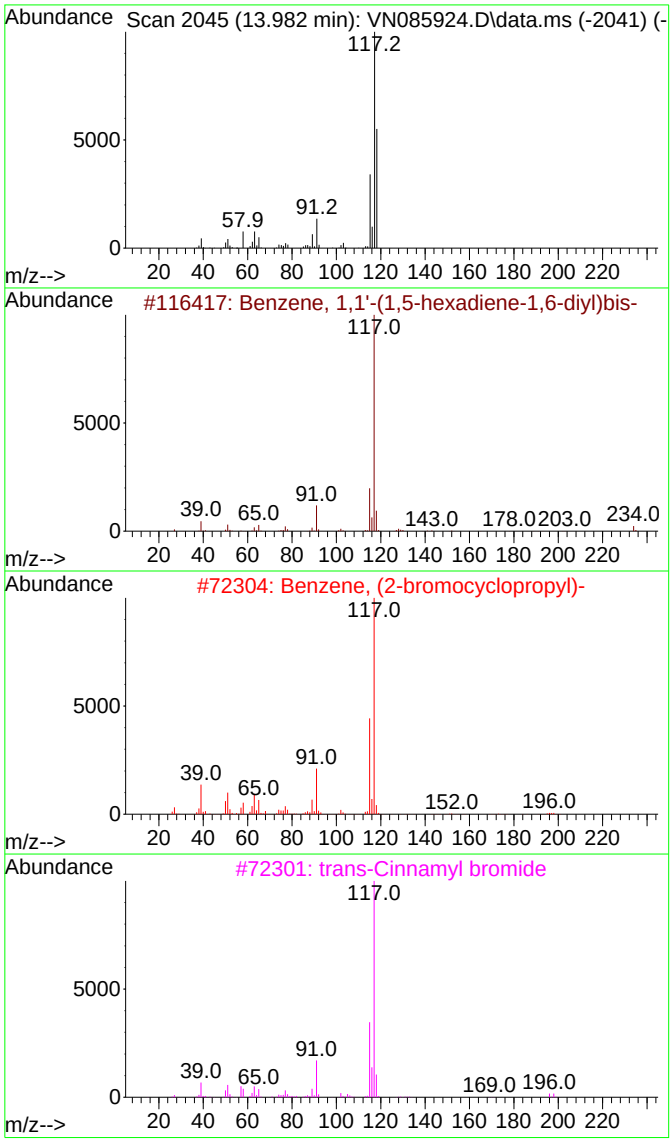
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

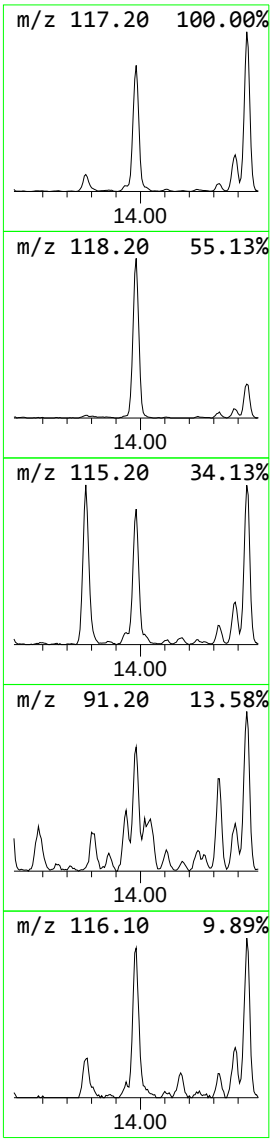
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.982	26.16 ug/l	635226	1,4-Dichlorobenzene-d4	13.776

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,1'-(1,5-hexadiene-1,6...	234	C18H18	004439-45-6	59
2	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	59
3	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	59
4	4-Ethynylaniline	117	C8H7N	014235-81-5	47
5	Indole	117	C8H7N	000120-72-9	43





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

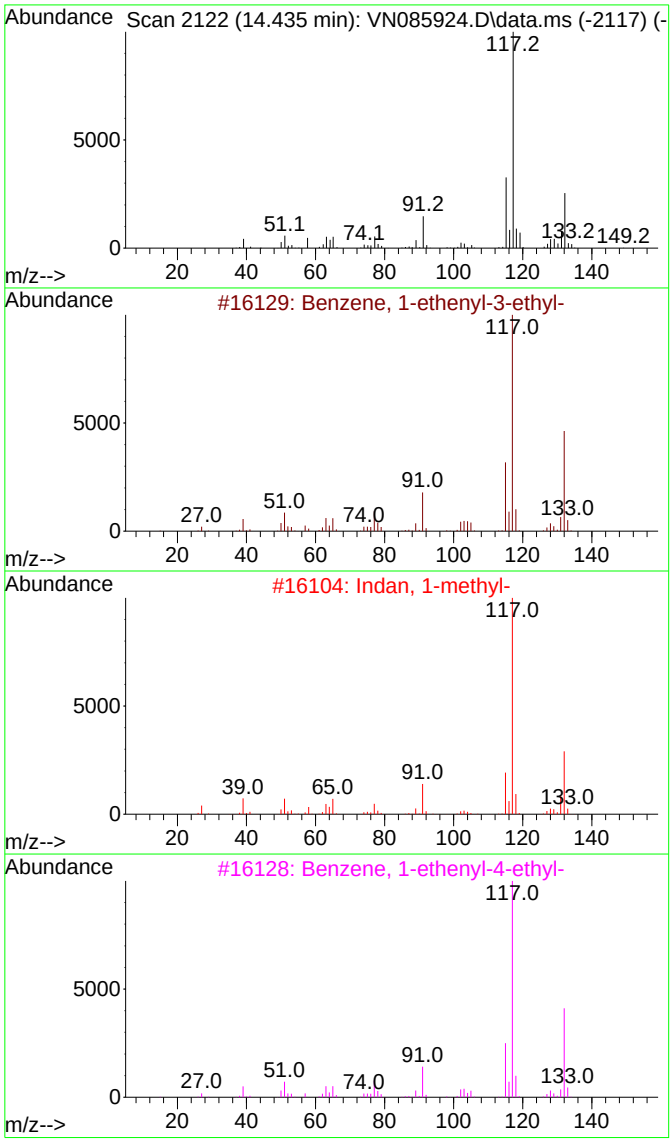
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

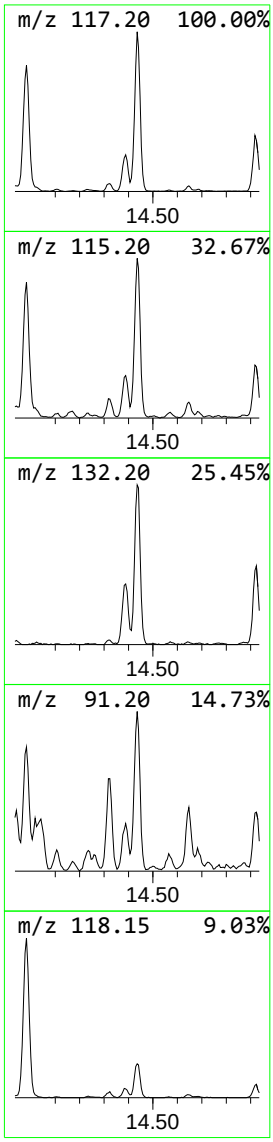
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.435	34.48 ug/l	837378	1,4-Dichlorobenzene-d4	13.776

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
2	Indan, 1-methyl-	132	C10H12	000767-58-8	87
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	86
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	83
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

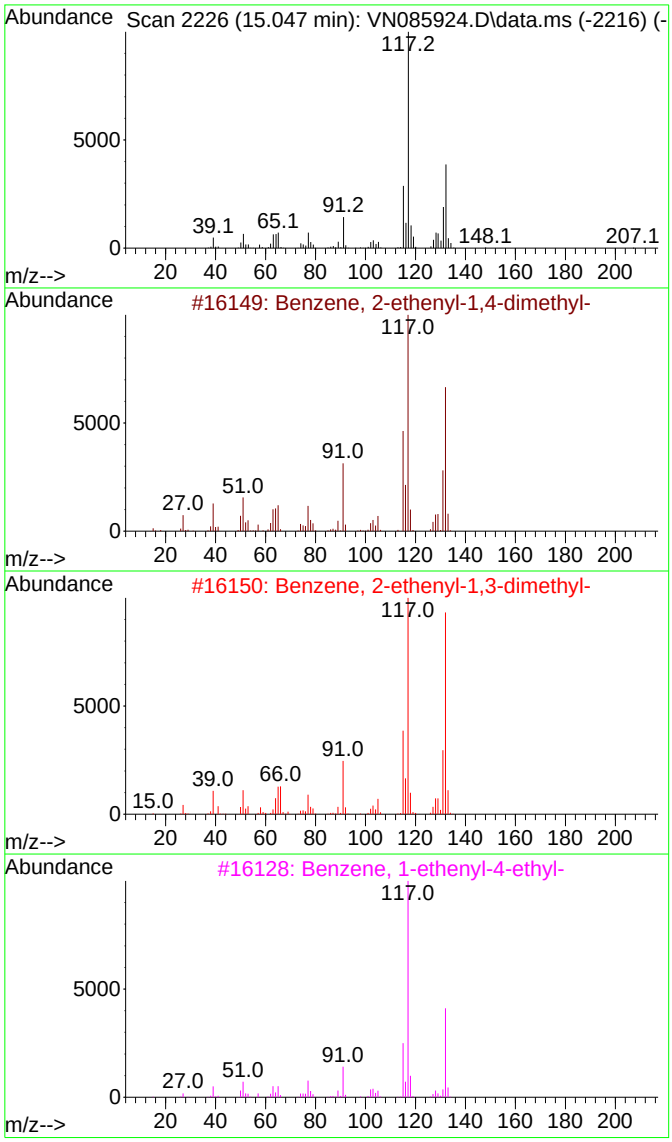
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.M
Quant Title : SW846 8260

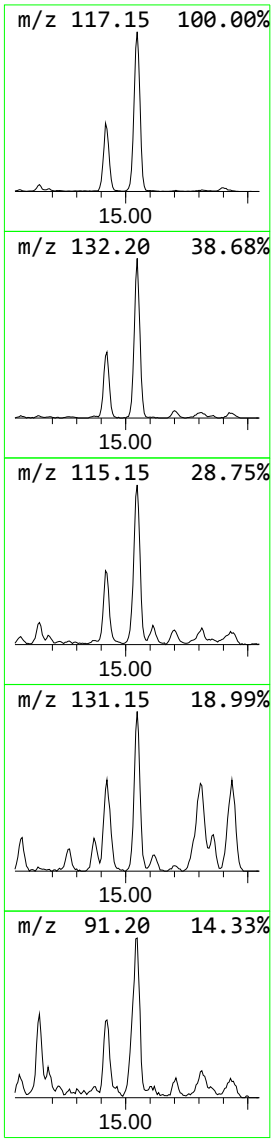
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.047	39.50 ug/l	959171	1,4-Dichlorobenzene-d4	13.776

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	94
3	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	91
5	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN031025\
Data File : VN085924.D
Acq On : 10 Mar 2025 19:12
Operator : JC\MD
Sample : Q1525-01
Misc : 5.0mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW10

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N021825W.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-	2.236	31.1	ug/l	559127	1	7.871	898542	50.0
Butane, 2,3-dim...	4.283	22.0	ug/l	395173	1	7.871	898542	50.0
Pentane, 2-methyl-	4.348	33.0	ug/l	593681	1	7.871	898542	50.0
Pentane, 3-methyl-	4.895	35.3	ug/l	634459	1	7.871	898542	50.0
Cyclopentane, m...	6.742	31.1	ug/l	558075	1	7.871	898542	50.0
Pentane, 2,3-di...	7.971	22.9	ug/l	411289	1	7.871	898542	50.0
Benzene, 1-ethy...	13.318	23.4	ug/l	568728	4	13.776	1214130	50.0
Benzene, 1,1'-(...	13.982	26.2	ug/l	635226	4	13.776	1214130	50.0
Benzene, 1-ethe...	14.435	34.5	ug/l	837378	4	13.776	1214130	50.0
Benzene, 2-ethe...	15.047	39.5	ug/l	959171	4	13.776	1214130	50.0