LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

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\82N122721W.M
Title : SW846 8260

Signal : TIC: VN070472.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	81	92	105	rBV5	126628	333540	6.31%	0.756%
2	2.446	161	170	189	rVB	162068	273899	5.18%	0.621%
3	3.140	411	429	456	rBV	815385	1892632	35.82%	4.290%
4	3.424	526	535	551	rVB3	29286	60691	1.15%	0.138%
5	3.518	553	570	602	rVB5	196351	597840	11.31%	1.355%
6	3.717	629	644	670	rVB2	84469	209764	3.97%	0.475%
7	3.891	693	709	728	rVB3	51010	134718	2.55%	0.305%
8	3.990	728	746	769	rVB3	44595	123186	2.33%	0.279%
9	4.259	818	846	881	rBV7	61668	274075	5.19%	0.621%
10	4.583	937	967	993	rBV2	30587	106156	2.01%	0.241%
11	4.953	1083	1105	1107	rBV8	26692	54482	1.03%	0.123%
12	5.079	1128	1152	1163	rBV3	133129	450028	8.52%	1.020%
13	5.618	1322	1353	1381	rBV2	161582	558353	10.57%	1.265%
14	5.935	1445	1471	1503	rBV4	36791	127114	2.41%	0.288%
15	6.120	1512	1540	1560	rBV4	45150	135873	2.57%	0.308%
16	6.289	1577	1603	1622	rBV7	26305	82144	1.55%	0.186%
17	6.401	1624	1645	1658	rBV4	27430	85041	1.61%	0.193%
18	6.605	1699	1721	1739	rBV4	55808	175299	3.32%	0.397%
19	6.696	1741	1755	1782	rVV6	31117	122327	2.32%	0.277%
20	6.900	1811	1831	1850	rBV6	34298	97790	1.85%	0.222%
21	7.045	1862	1885	1902	rBV2	89462	259213	4.91%	0.587%
22	7.144	1905	1922	1943	rVV3	120859	337433	6.39%	0.765%
23	7.316	1972	1986	2011	rVB5	20217	62025	1.17%	0.141%
24	7.844	2171	2183	2202	rVB5	26812	65831	1.25%	0.149%
25	8.021	2224	2249	2259	rBV	640392	1557301	29.47%	3.530%
26	8.083	2260	2272	2292	rVV2	1363026	3259324	61.68%	7.387%
27	8.172	2293	2305	2333	rVB4	188414	491756	9.31%	1.115%
28	8.295	2333	2351	2367	rBV4	113678	261659	4.95%	0.593%
29	8.437	2384	2404	2428	rBV	817842	1922235	36.38%	4.357%
30	8.614	2438	2470	2494	rVB	645711	1825081	34.54%	4.136%
31	8.826	2530	2549	2572	rVB7	35233	98687	1.87%	0.224%
32	8.965	2580	2601	2629	rBV	2049091	4352339	82.37%	9.864%
33	9.244	2697	2705	2727	rVB4	36128	78323	1.48%	0.178%
34	9.456	2754	2784	2795	rBV5	119489	335898	6.36%	0.761%
35	9.510	2795	2804	2821	rVV4	81555	164710	3.12%	0.373%
36	9.730	2866	2886	2899	rBV6	22677	55189	1.04%	0.125%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
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Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

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\82N122721W.M
Title : SW846 8260

37	9.880	2928	2942	2965	rVB	257593	521846	9.88%	1.183%
38	10.014	2966	2992	3006	rBV3	278076	610088	11.55%	1.383%
39	10.078	3007	3016	3024	rBV3	39630	64580	1.22%	0.146%
40	10.213	3047	3066	3092	rBV5	78425	218223	4.13%	0.495%
41	10.368	3114	3124	3134	rVB2	58463	95759	1.81%	0.217%
42	10.440	3134	3151	3171	rBV	2797415	5284092	100.00%	11.976%
43	10.620	3206	3218	3232	rVB9	39736	95612	1.81%	0.217%
44	11.524	3537	3555	3570	rBV5	30076	68851	1.30%	0.156%
45	11.623	3582	3592	3618	rVB3	28763	67868	1.28%	0.154%
46	11.741	3618	3636	3662	rBV	2573792	4745600	89.81%	10.756%
47	11.843	3663	3674	3695	rVV	131321	254237	4.81%	0.576%
48	11.953	3698	3715	3739	rVB2	262974	569898	10.79%	1.292%
49	12.278	3812	3836	3857	rBV	159067	305930	5.79%	0.693%
50	12.575	3934	3947	3961	rBV2	141126	240138	4.54%	0.544%
51	12.728	3989	4004	4034	rVB	1903102	3444541	65.19%	7.807%
52	12.919	4063	4075	4087	rVV	62131	101952	1.93%	0.231%
53	12.980	4087	4098	4103	rVV2	44573	70912	1.34%	0.161%
54	13.013	4104	4110	4118	rVV3	77860	128379	2.43%	0.291%
55	13.055	4119	4126	4143	rVB2	63684	111067	2.10%	0.252%
56	13.219	4171	4187	4205	rBV	318855	554377	10.49%	1.256%
57	13.367	4228	4242	4259	rVB	437065	748697	14.17%	1.697%
58	13.672	4340	4356	4384	rBV	1938505	3625971	68.62%	8.218%
59	13.873	4422	4431	4440	rVV	75212	124236	2.35%	0.282%
60	14.061	4486	4501	4513	rBV3	44393	80373	1.52%	0.182%
61	14.128	4514	4526	4532	rBV2	51200	87695	1.66%	0.199%
62	14.214	4548	4558	4573	rVB2	74619	118563	2.24%	0.269%
63	14.324	4587	4599	4615	rVB2	102382	180539	3.42%	0.409%
64	14.536	4664	4678	4685	rBV2	39178	67646	1.28%	0.153%
65	14.576	4685	4693	4702	rVB2	47232	74602	1.41%	0.169%
66	14.809	4769	4780	4791	rBV3	39305	68150	1.29%	0.154%
67	14.927	4806	4824	4837	rBV4	68128	160495	3.04%	0.364%
68	15.193	4901	4923	4948	rBV8	24122	78886	1.49%	0.179%
69	15.507	5025	5040	5056	rVB	67269	129883	2.46%	0.294%

Sum of corrected areas: 44121642

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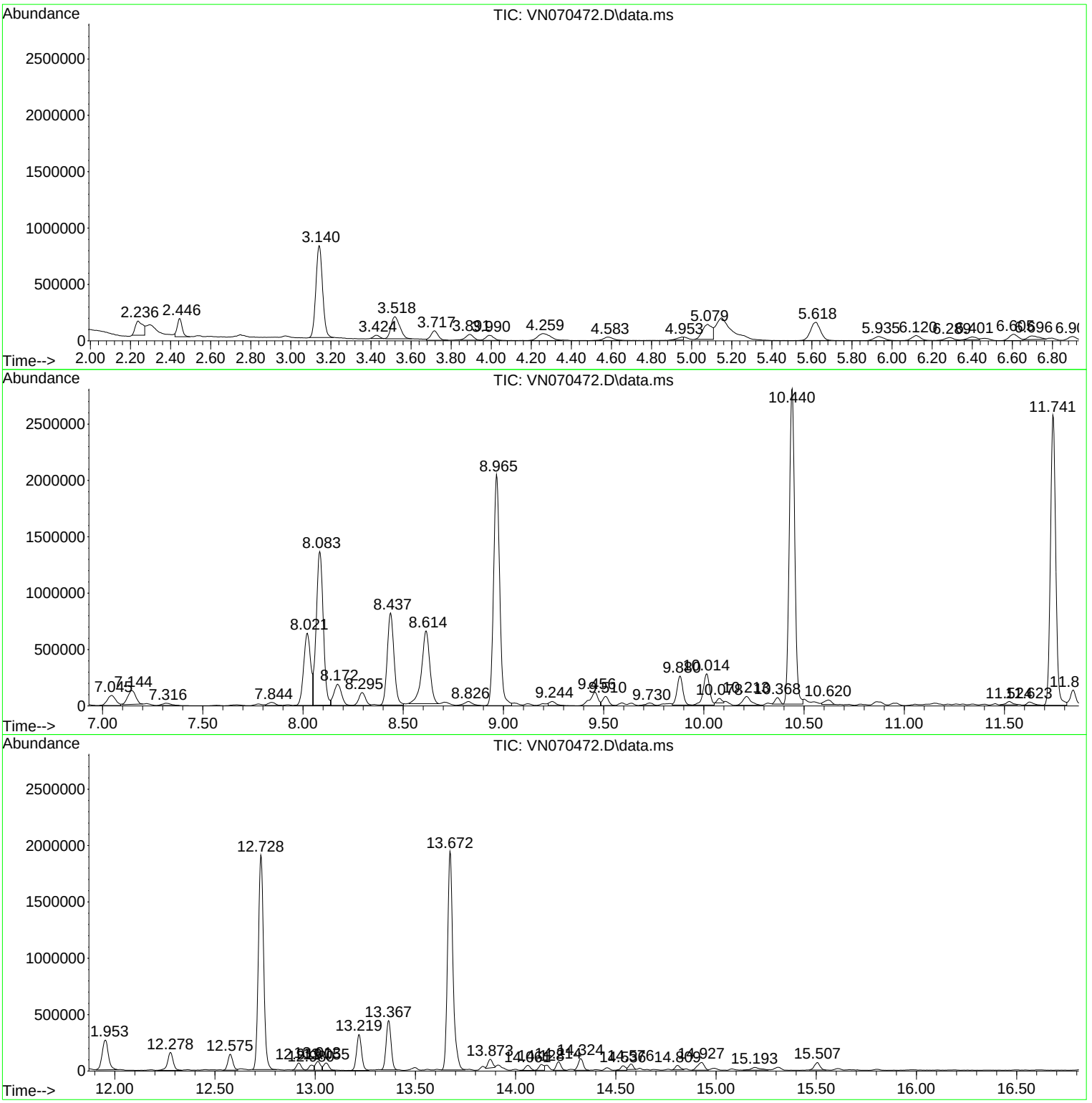
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
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ClientSampleId :
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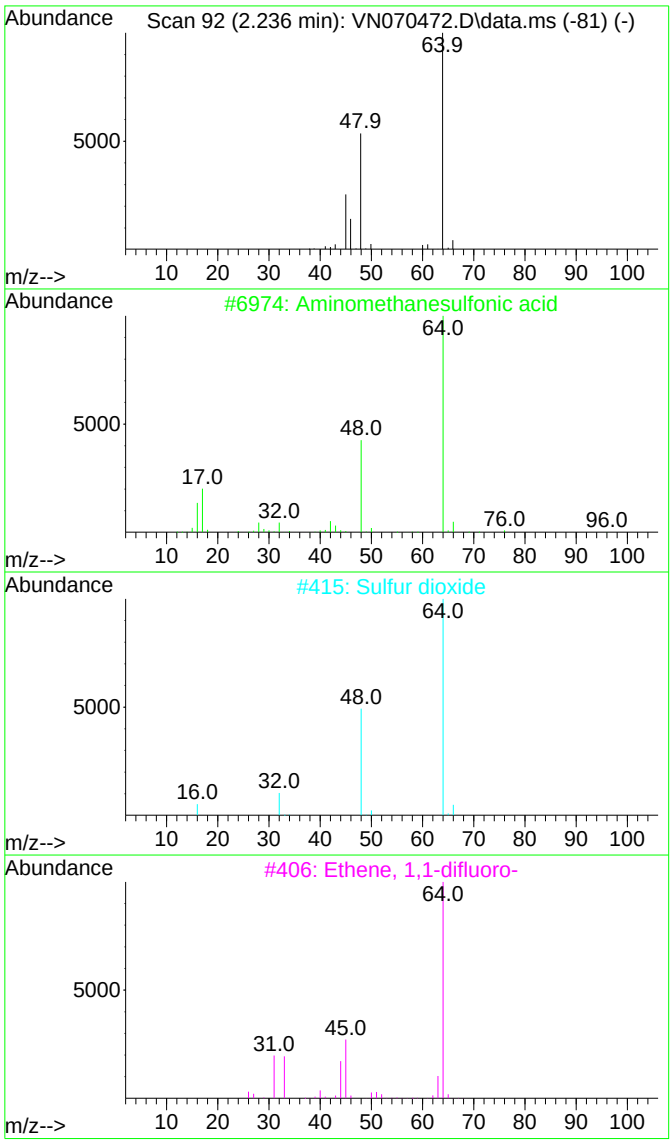
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

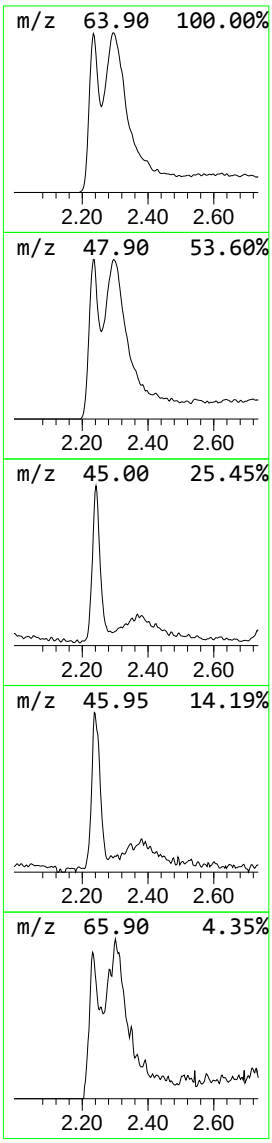
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Aminomethanesulfonic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.236	5.12 ug/l	333540	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	78
2	Sulfur dioxide	64	O2S	007446-09-5	78
3	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	9
4	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	7





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

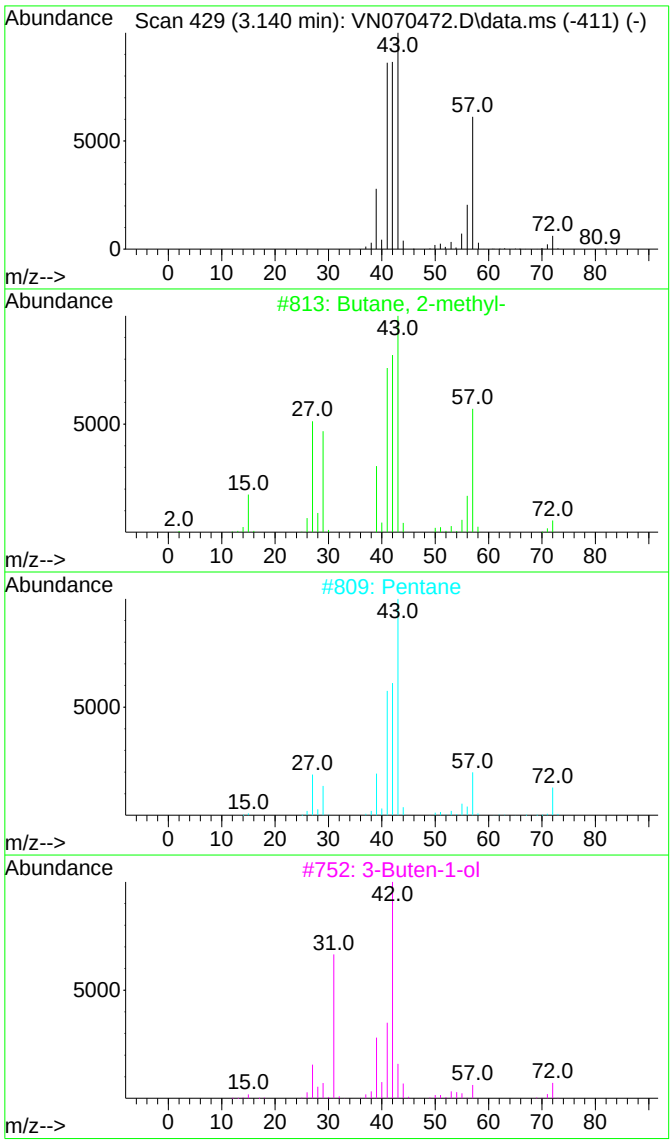
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

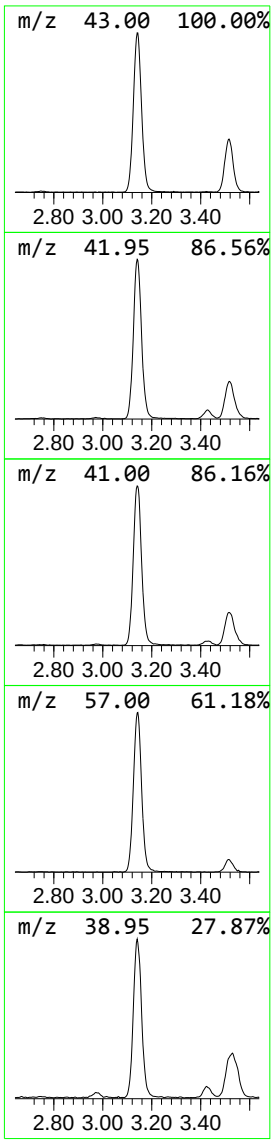
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.140	29.03 ug/l	1892630	Pentafluorobenzene	8.083

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	50
3	3-Buten-1-ol	72	C4H8O	000627-27-0	16
4	1-Butene	56	C4H8	000106-98-9	9
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

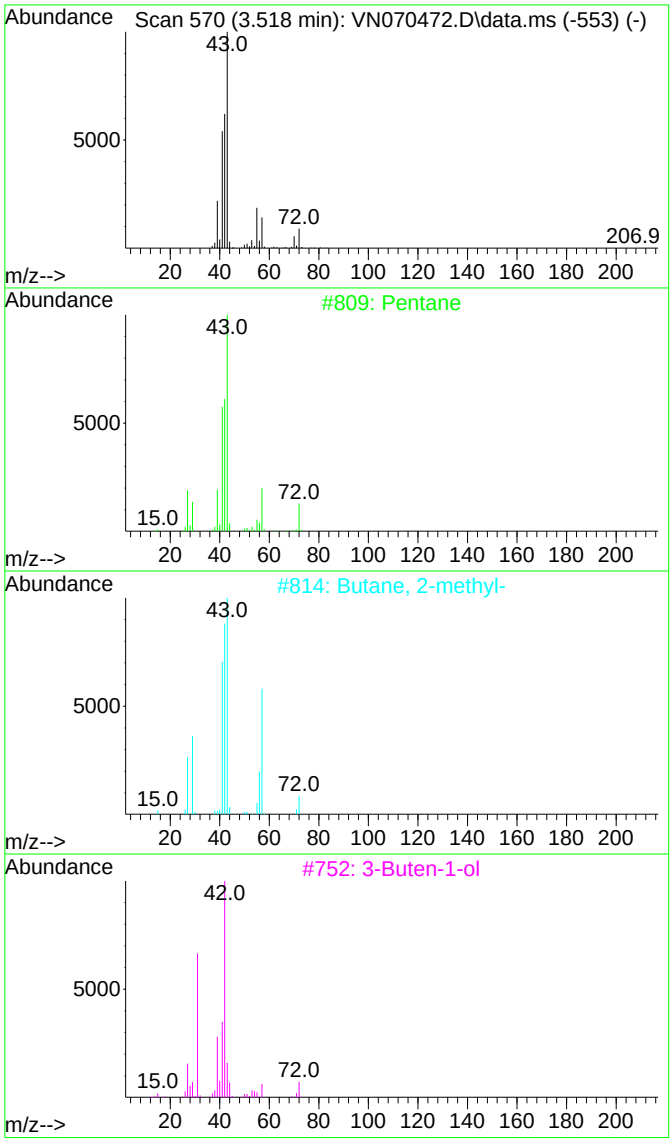
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

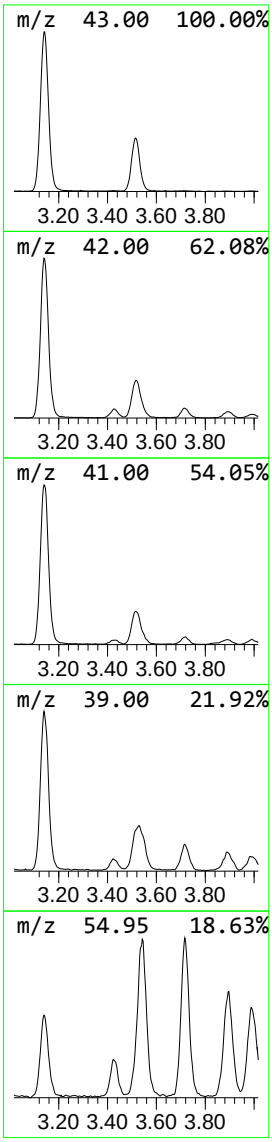
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.518	9.17 ug/l	597840	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	80
2	Butane, 2-methyl-	72	C5H12	000078-78-4	53
3	3-Buten-1-ol	72	C4H8O	000627-27-0	47
4	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	32
5	Oxirane, ethyl-	72	C4H8O	000106-88-7	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

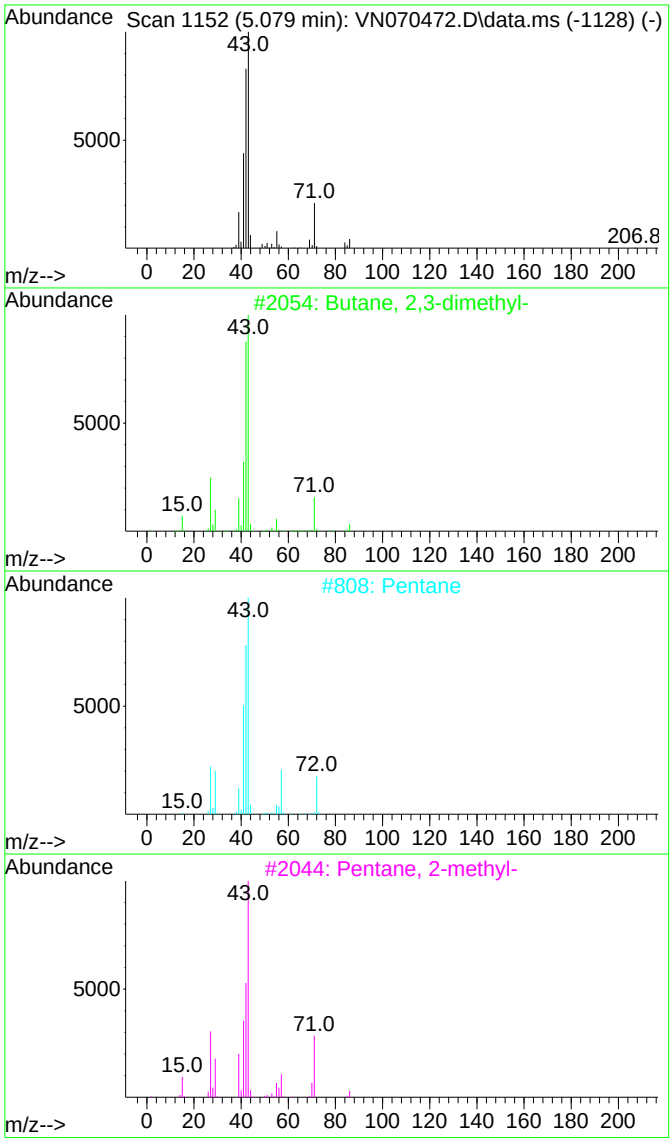
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

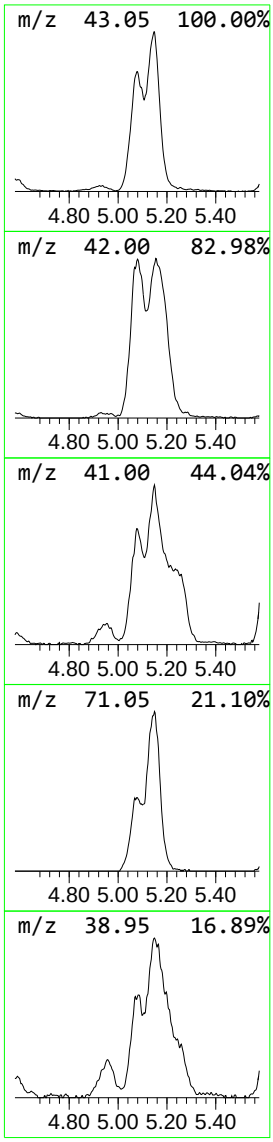
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Butane, 2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.079	6.90 ug/l	450028	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	86
2	Pentane	72	C5H12	000109-66-0	42
3	Pentane, 2-methyl-	86	C6H14	000107-83-5	38
4	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9
5	Hydroxylamine, O-(3-methylbutyl)-	103	C5H13NO	019411-65-5	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

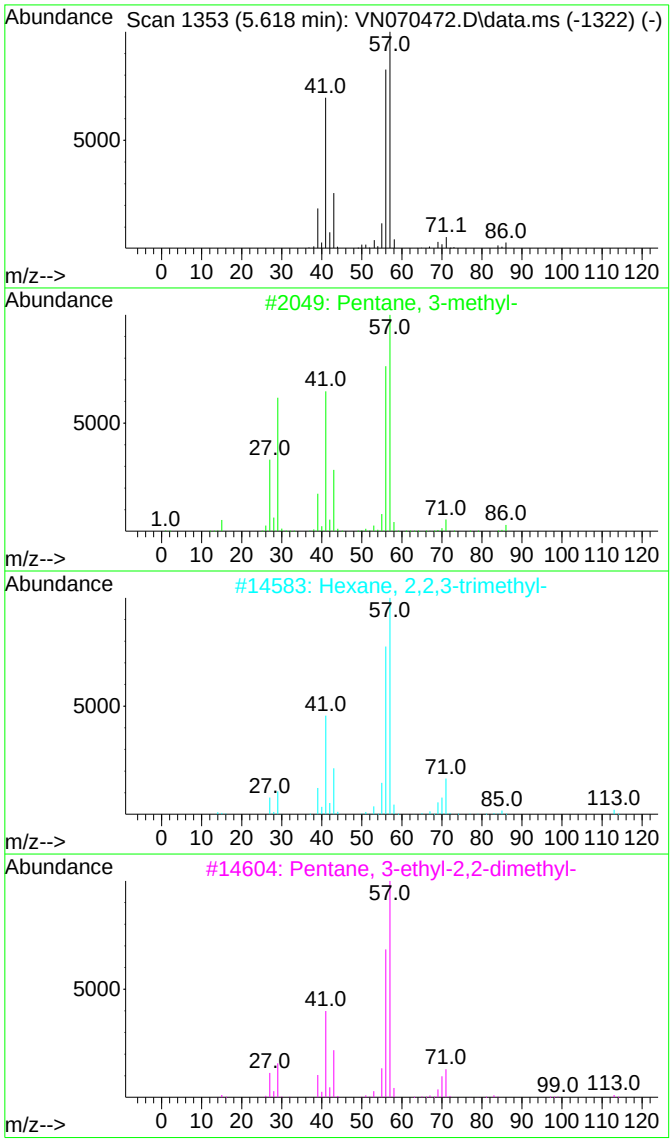
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

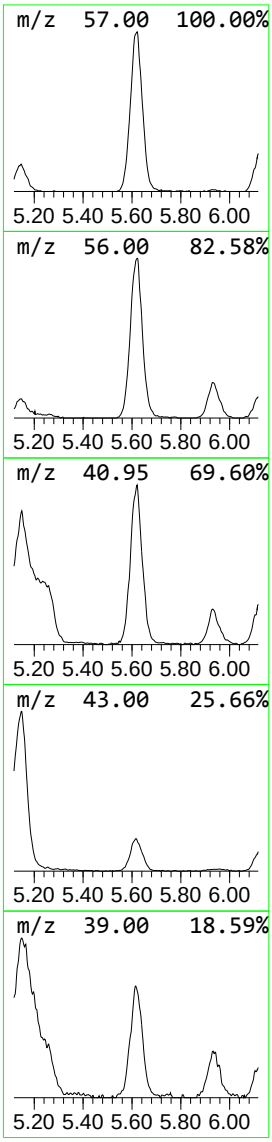
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.618	8.57 ug/l	558353	Pentafluorobenzene	8.083

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	78
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	45
5	Butyl isocyanatoacetate	157	C7H11NO3	017046-22-9	43





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ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

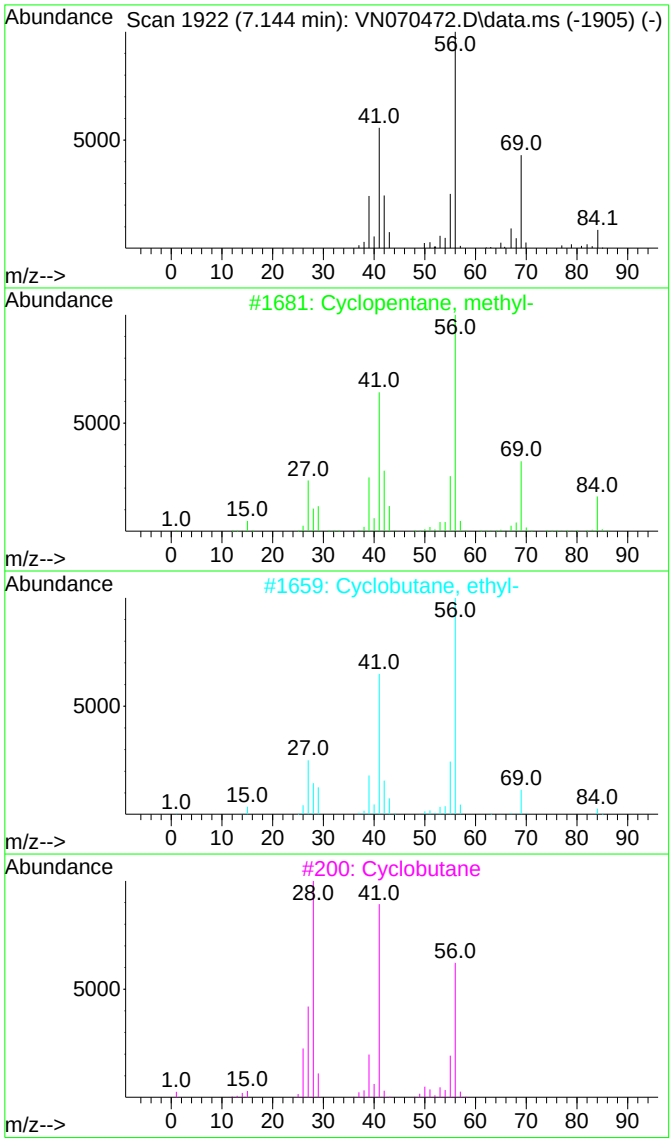
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

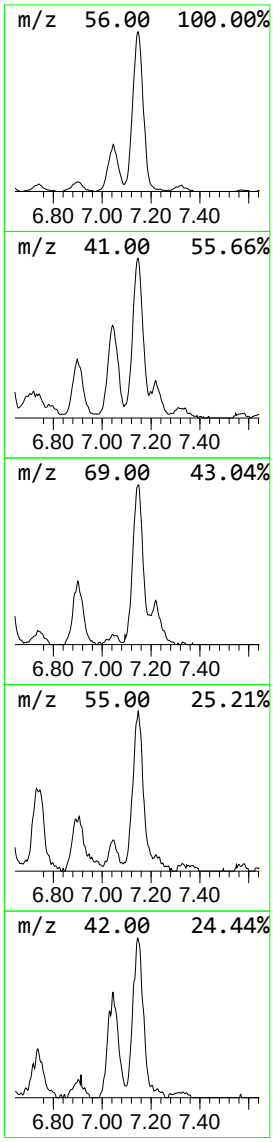
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.144	5.18 ug/l	337433	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	86
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
3	Cyclobutane	56	C4H8	000287-23-0	53
4	Cyclobutane, butyl-	112	C8H16	013152-44-8	45
5	Cyclobutanecarbonitrile, 3,3-dim...	109	C7H11N	053783-86-1	40





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

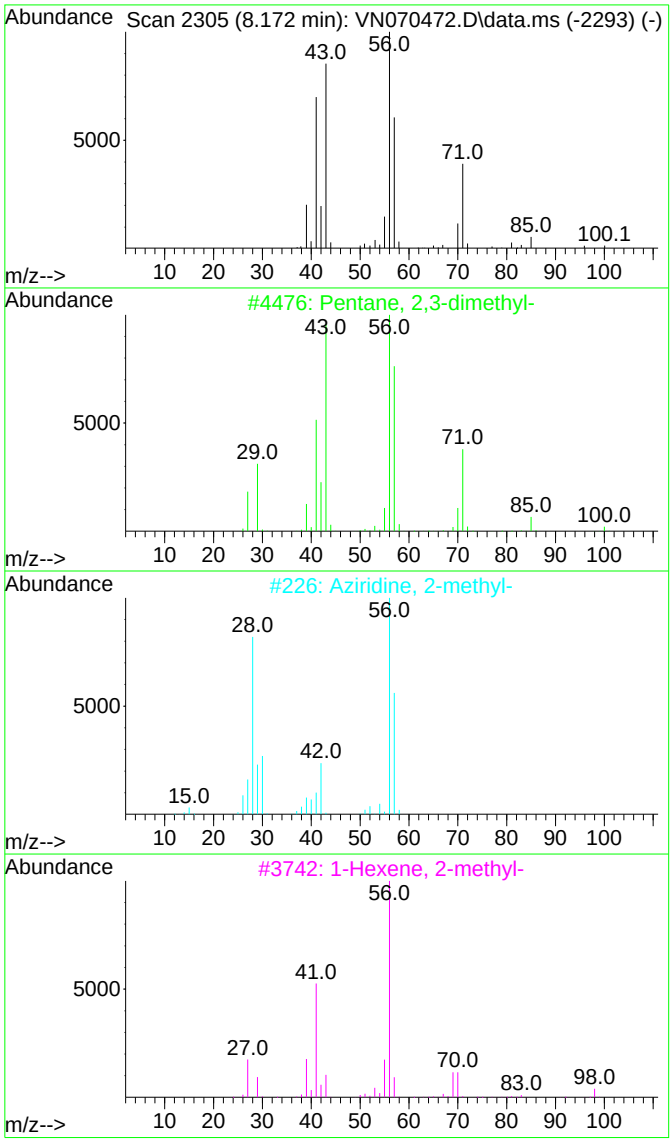
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

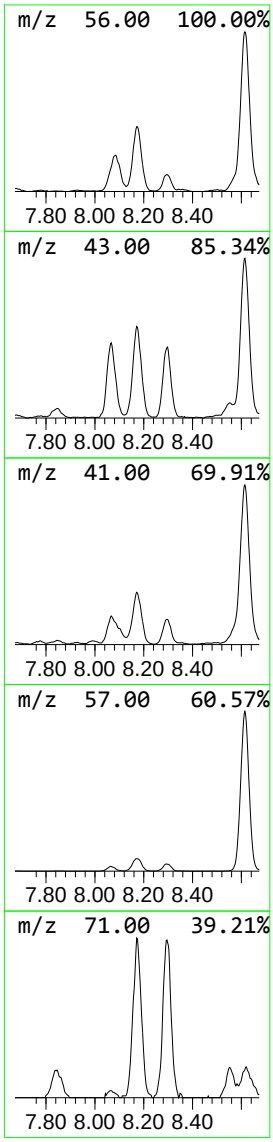
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentane, 2,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.172	7.54 ug/l	491756	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	43
3	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	37
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

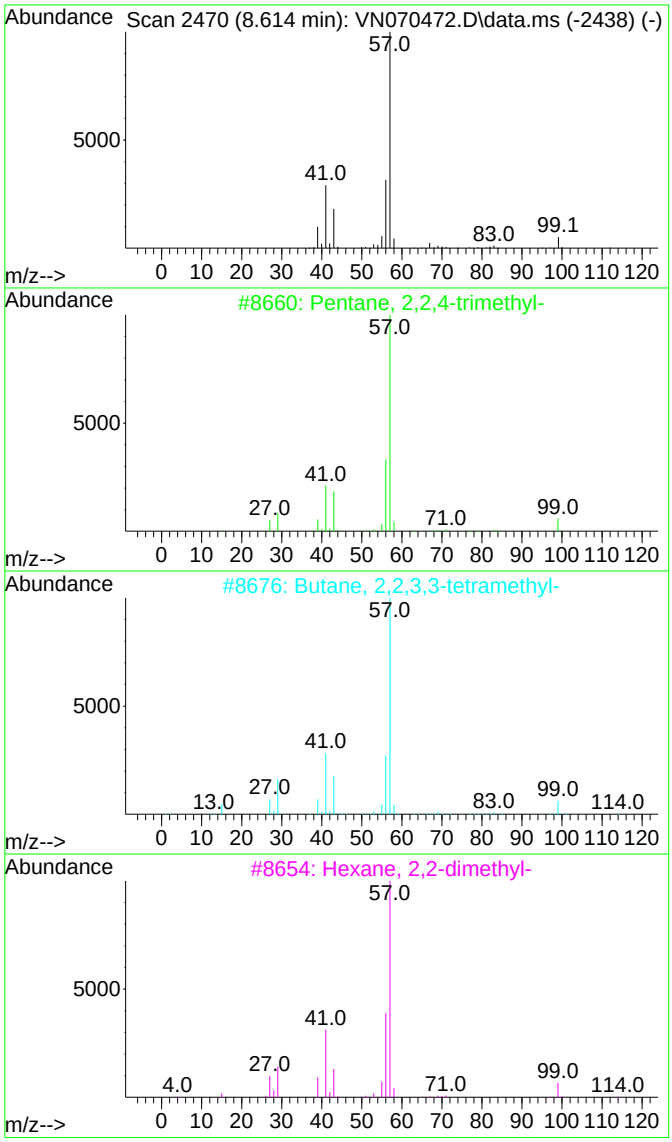
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

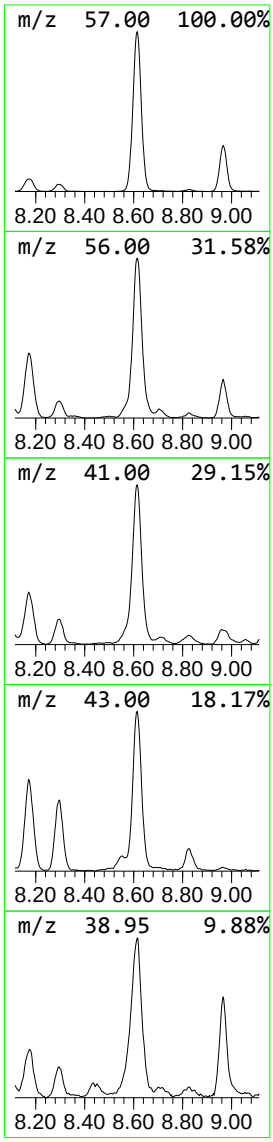
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Pentane, 2,2,4-trimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.614	20.97 ug/l	1825080	1,4-Difluorobenzene	8.965

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
2	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	72
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

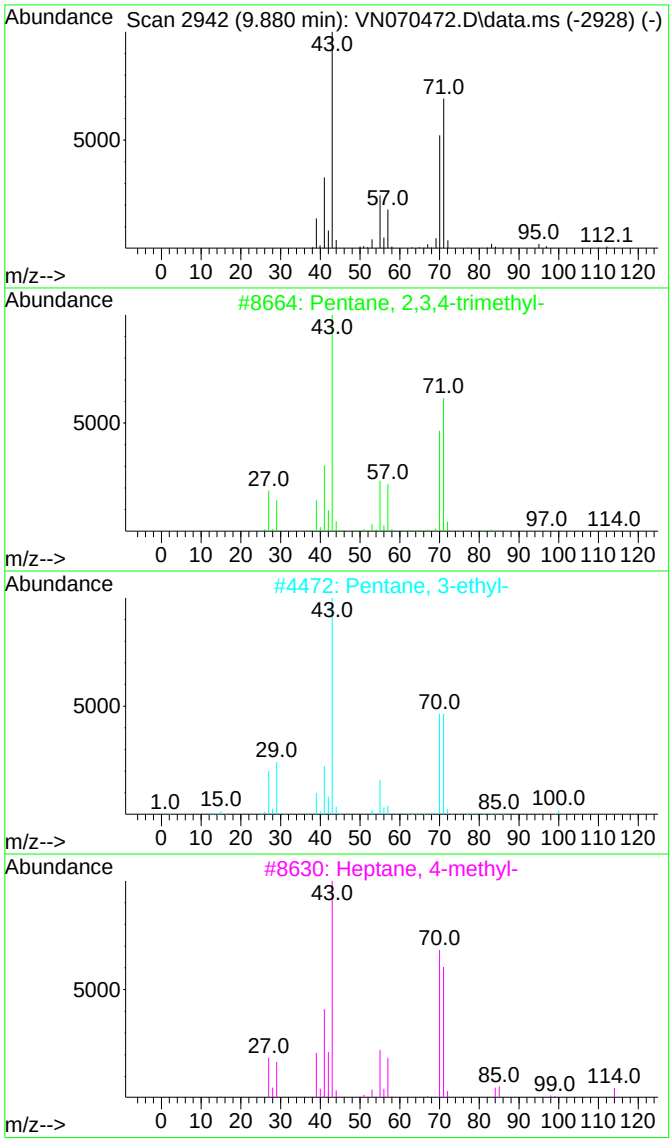
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

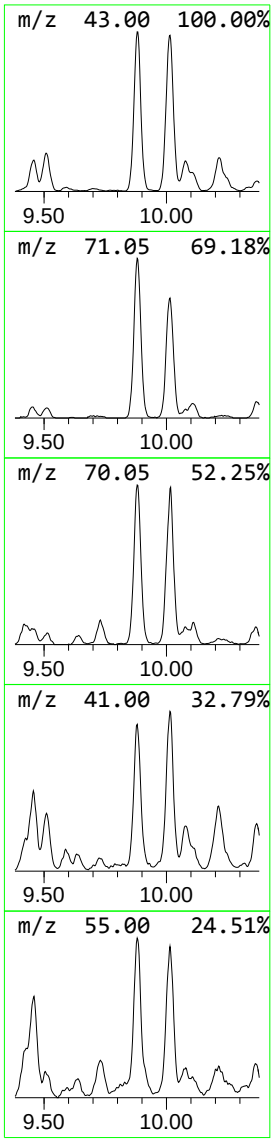
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Pentane, 2,3,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.880	6.00 ug/l	521846	1,4-Difluorobenzene	8.965

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	90
2	Pentane, 3-ethyl-	100	C7H16	000617-78-7	72
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	64
5	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

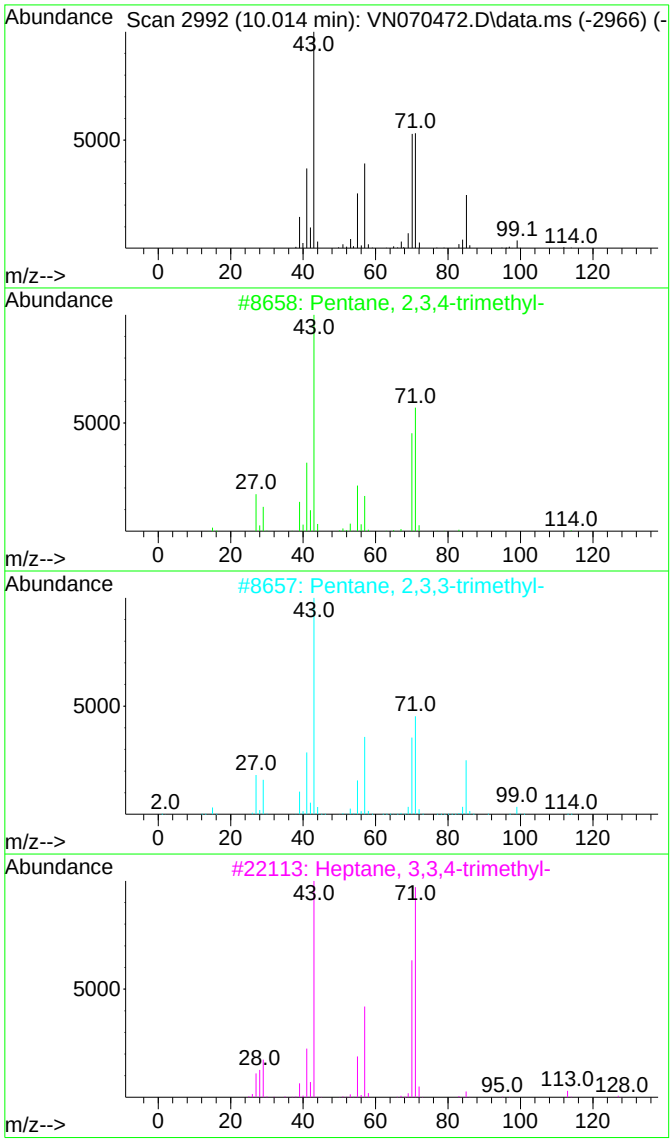
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

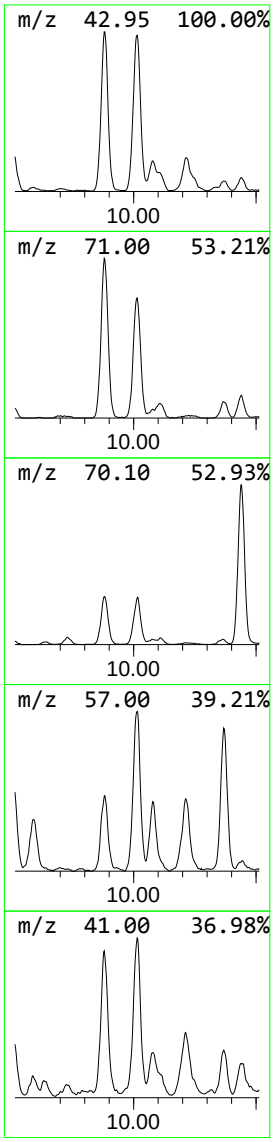
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Pentane, 2,3,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.014	7.01 ug/l	610088	1,4-Difluorobenzene	8.965

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	74
2	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
3	Heptane, 3,3,4-trimethyl-	142	C10H22	020278-87-9	59
4	1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	59
5	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

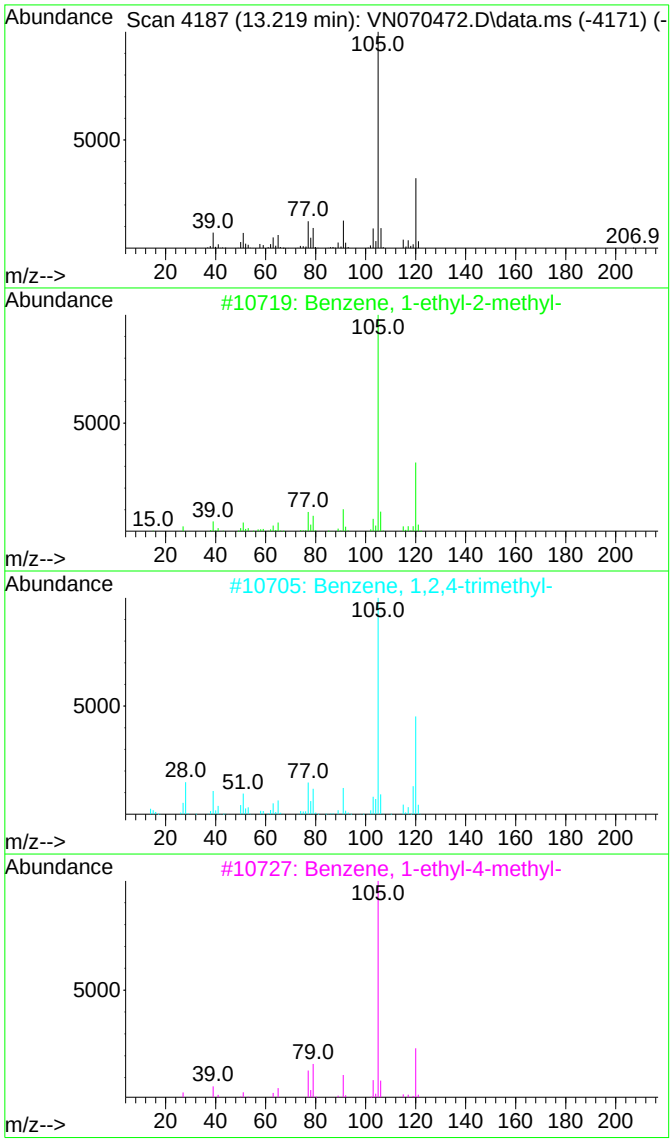
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

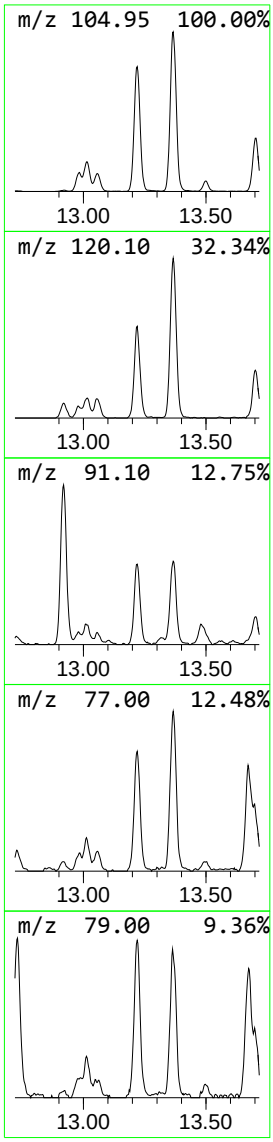
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.219	7.64 ug/l	554377	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010422\
Data File : VN070472.D
Acq On : 4 Jan 2022 18:49
Operator : JC/MD
Sample : M5251-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 21 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.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
Aminomethanesul...	2.236	5.1	ug/l	333540	1	8.083	3259320	50.0
Butane, 2-methyl-	3.140	29.0	ug/l	1892630	1	8.083	3259320	50.0
Pentane	3.518	9.2	ug/l	597840	1	8.083	3259320	50.0
Butane, 2,3-dim...	5.079	6.9	ug/l	450028	1	8.083	3259320	50.0
Pentane, 3-methyl-	5.618	8.6	ug/l	558353	1	8.083	3259320	50.0
Cyclopentane, m...	7.144	5.2	ug/l	337433	1	8.083	3259320	50.0
Pentane, 2,3-di...	8.172	7.5	ug/l	491756	1	8.083	3259320	50.0
Pentane, 2,2,4-...	8.614	21.0	ug/l	1825080	2	8.965	4352340	50.0
Pentane, 2,3,4-...	9.880	6.0	ug/l	521846	2	8.965	4352340	50.0
Pentane, 2,3,3-...	10.014	7.0	ug/l	610088	2	8.965	4352340	50.0
Benzene, 1-ethy...	13.219	7.6	ug/l	554377	4	13.672	3625970	50.0