LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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: VN070434.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.258	83	100	111	rBV	449372	883483	8.18%	0.668%
2	2.443	160	169	191	rVB	382944	634981	5.88%	0.480%
3	3.138	407	428	471	rBV	4612032	10806377	100.00%	8.165%
4	3.510	555	567	590	rVB2	109389	267846	2.48%	0.202%
5	3.843	670	691	719	rVB3	45152	148687	1.38%	0.112%
6	3.988	721	745	773	rBV4	89175	249523	2.31%	0.189%
7	4.251	810	843	883	rBV	842054	2984947	27.62%	2.255%
8	4.580	946	966	990	rBV2	58967	184038	1.70%	0.139%
9	5.077	1118	1151	1174	rBV2	1862424	7548188	69.85%	5.703%
10	5.616	1313	1352	1392	rBV2	639568	2223707	20.58%	1.680%
11	6.605	1699	1721	1746	rBV3	115639	360060	3.33%	0.272%
12	6.890	1803	1827	1849	rBV6	157936	537144	4.97%	0.406%
13	7.045	1862	1885	1903	rBV2	616880	1840969	17.04%	1.391%
14	7.144	1904	1922	1938	rVV	723998	2139757	19.80%	1.617%
15	7.219	1940	1950	1972	rVV3	398260	1062202	9.83%	0.803%
16	7.324	1975	1989	2022	rVB8	74396	266632	2.47%	0.201%
17	7.651	2094	2111	2141	rVB5	57757	179157	1.66%	0.135%
18	7.842	2168	2182	2208	rVB4	323570	882817	8.17%	0.667%
19	8.019	2220	2248	2259	rBV	744139	1846678	17.09%	1.395%
20	8.086	2260	2273	2289	rVV3	1605053	3887549	35.97%	2.937%
21	8.172	2289	2305	2333	rVV2	1765157	4547215	42.08%	3.436%
22	8.295	2336	2351	2361	rVV4	78721	186717	1.73%	0.141%
23	8.357	2362	2374	2386	rVV4	111911	270035	2.50%	0.204%
24	8.456	2387	2411	2433	rVV4	2756718	7715791	71.40%	5.830%
25	8.614	2434	2470	2491	rVV	2082406	6248806	57.83%	4.722%
26	8.705	2492	2504	2526	rVB3	433263	997472	9.23%	0.754%
27	8.965	2581	2601	2626	rBV	2214449	5046492	46.70%	3.813%
28	9.059	2628	2636	2651	rVB2	195334	391110	3.62%	0.296%
29	9.196	2672	2687	2696	rBV2	267000	550158	5.09%	0.416%
30	9.242	2698	2704	2724	rVB2	226237	445614	4.12%	0.337%
31	9.421	2750	2771	2781	rBV2	671355	1573641	14.56%	1.189%
32	9.513	2796	2805	2822	rVB5	218067	440887	4.08%	0.333%
33	9.590	2822	2834	2844	rBV4	90105	174256	1.61%	0.132%
34	9.732	2866	2887	2901	rBV4	180529	439867	4.07%	0.332%
35	9.826	2901	2922	2928	rBV8	82418	243364	2.25%	0.184%
36	9.880	2929	2942	2961	rVB	938282	1913047	17.70%	1.446%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
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Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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	9.974	2962	2977	2980	rBV4	305991	482609	4.47%	0.365%
38	10.011	2982	2991	3020	rVB	965484	2054893	19.02%	1.553%
39	10.119	3021	3031	3045	rVB8	80648	190035	1.76%	0.144%
40	10.196	3046	3060	3074	rBV6	193133	383322	3.55%	0.290%
41	10.320	3093	3106	3115	rBV	386806	709364	6.56%	0.536%
42	10.441	3135	3151	3195	rBV	3331221	6702839	62.03%	5.065%
43	10.639	3200	3225	3240	rVV2	1096709	2350932	21.76%	1.776%
44	10.714	3240	3253	3268	rVV3	965879	1748958	16.18%	1.322%
45	10.778	3268	3277	3290	rVV2	96606	169673	1.57%	0.128%
46	10.856	3291	3306	3327	rVB5	705275	1618733	14.98%	1.223%
47	10.942	3328	3338	3358	rVB9	90735	228529	2.11%	0.173%
48	11.127	3395	3407	3417	rBV5	147087	273096	2.53%	0.206%
49	11.256	3446	3455	3467	rVB5	149944	255489	2.36%	0.193%
50	11.449	3517	3527	3541	rVB3	131931	229620	2.12%	0.174%
51	11.554	3556	3566	3586	rVB3	49685	129591	1.20%	0.098%
52	11.741	3621	3636	3652	rBV	2778472	5038237	46.62%	3.807%
53	11.843	3660	3674	3697	rVB	3291296	5828365	53.93%	4.404%
54	12.575	3933	3947	3967	rBV	1807903	3058865	28.31%	2.311%
55	12.728	3990	4004	4026	rVB	2357431	4159163	38.49%	3.143%
56	12.919	4061	4075	4092	rVB	1242970	2105129	19.48%	1.591%
57	13.219	4174	4187	4203	rBV	318485	543286	5.03%	0.411%
58	13.367	4233	4242	4255	rVB	117872	213039	1.97%	0.161%
59	13.482	4272	4285	4304	rVB3	317941	746087	6.90%	0.564%
60	13.672	4341	4356	4368	rBV	2174637	3867605	35.79%	2.922%
61	13.769	4384	4392	4403	rVB	144886	227812	2.11%	0.172%
62	13.836	4404	4417	4421	rBV	799695	1162799	10.76%	0.879%
63	13.873	4422	4431	4444	rVB	3947109	6509168	60.23%	4.918%
64	13.999	4466	4478	4491	rVB	643009	1030494	9.54%	0.779%
65	14.067	4491	4503	4517	rVB	446000	734560	6.80%	0.555%
66	14.147	4517	4533	4546	rBV3	390680	745193	6.90%	0.563%
67	14.214	4546	4558	4570	rBV2	348948	560636	5.19%	0.424%
68	14.278	4570	4582	4588	rBV2	207707	335945	3.11%	0.254%
69	14.327	4588	4600	4614	rVB	1906265	3131089	28.97%	2.366%
70	14.404	4618	4629	4640	rVB5	195544	358307	3.32%	0.271%
71	14.463	4641	4651	4661	rVB	169938	247660	2.29%	0.187%
72	14.536	4662	4678	4689	rBV	578128	974523	9.02%	0.736%
73	14.654	4715	4722	4733	rVB2	97771	143634	1.33%	0.109%
74	14.761	4750	4762	4770	rBV2	177663	282515	2.61%	0.213%
75	14.809	4770	4780	4806	rVB4	344728	723147	6.69%	0.546%
76	14.927	4807	4824	4838	rBV4	306701	699984	6.48%	0.529%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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

77	14.997	4838	4850	4865	rVB5	138756	244651	2.26%	0.185%
78	15.080	4871	4881	4891	rBV4	105562	185017	1.71%	0.140%
79	15.193	4910	4923	4934	rBV4	205231	389766	3.61%	0.295%
80	15.560	5050	5060	5072	rVV4	63753	108870	1.01%	0.082%
81	15.802	5136	5150	5163	rBV3	67133	135719	1.26%	0.103%
82	16.108	5252	5264	5278	rBV2	95871	206292	1.91%	0.156%

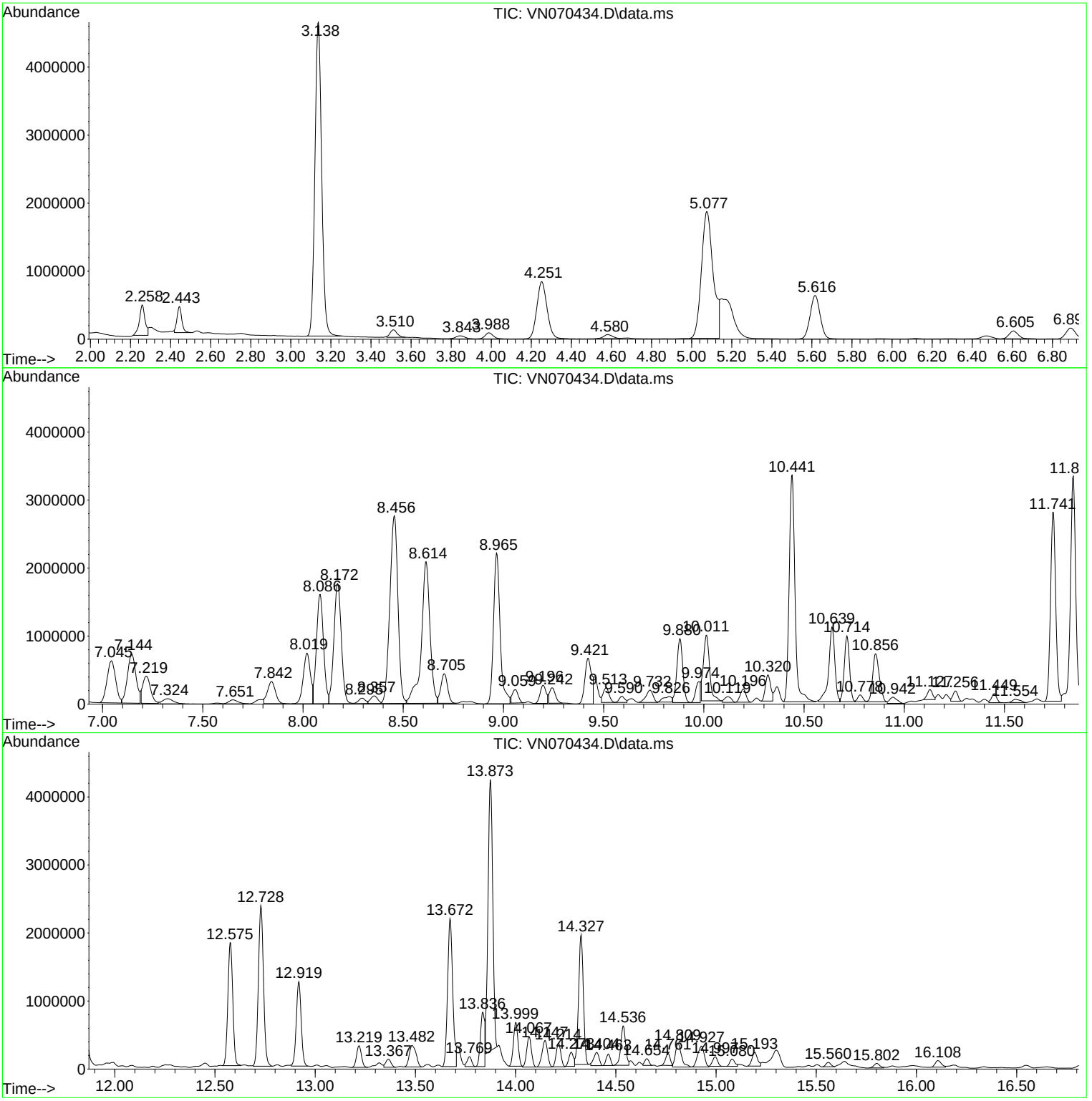
Sum of corrected areas: 132344454

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

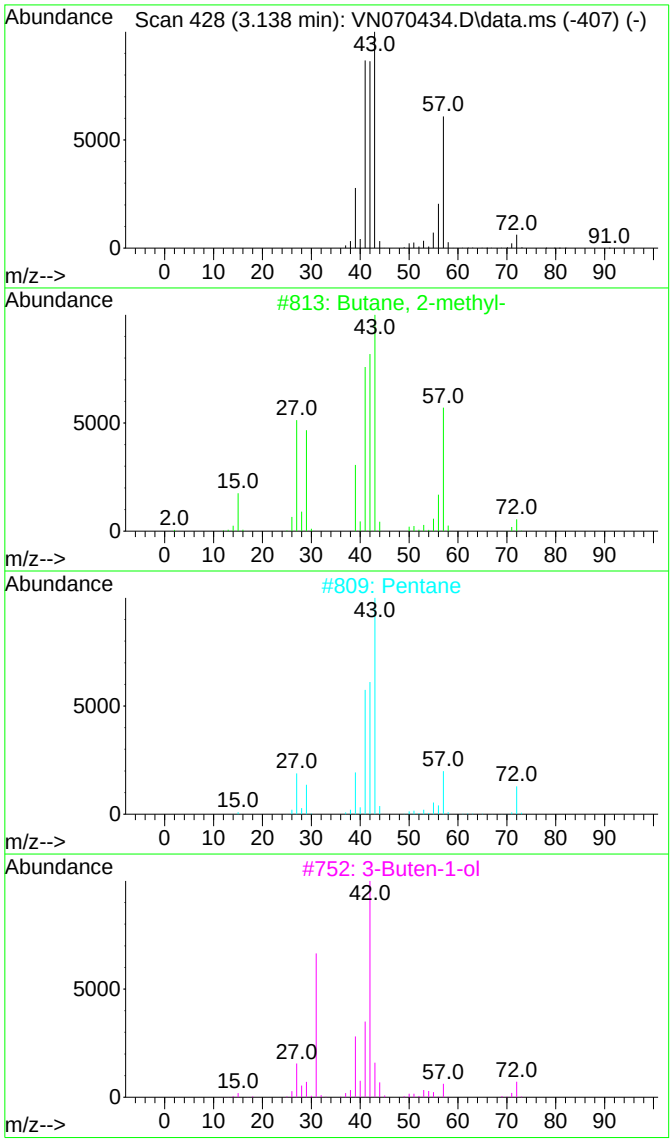
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

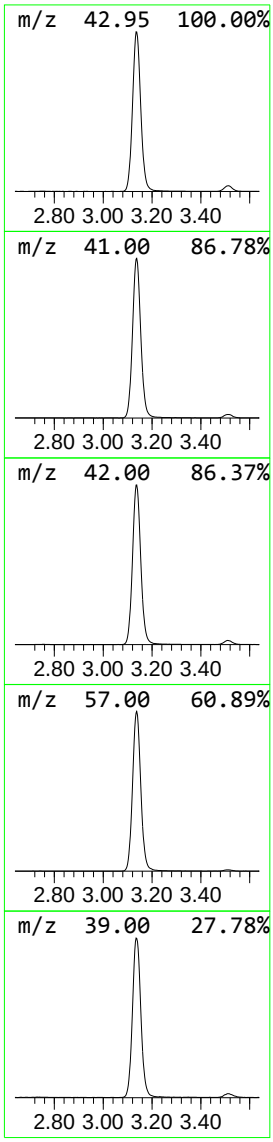
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.138	138.99 ug/l	10806400	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2	Pentane	72	C5H12	000109-66-0	36
3	3-Buten-1-ol	72	C4H8O	000627-27-0	28
4	Methane, isocyanato-	57	C2H3NO	000624-83-9	9
5	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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

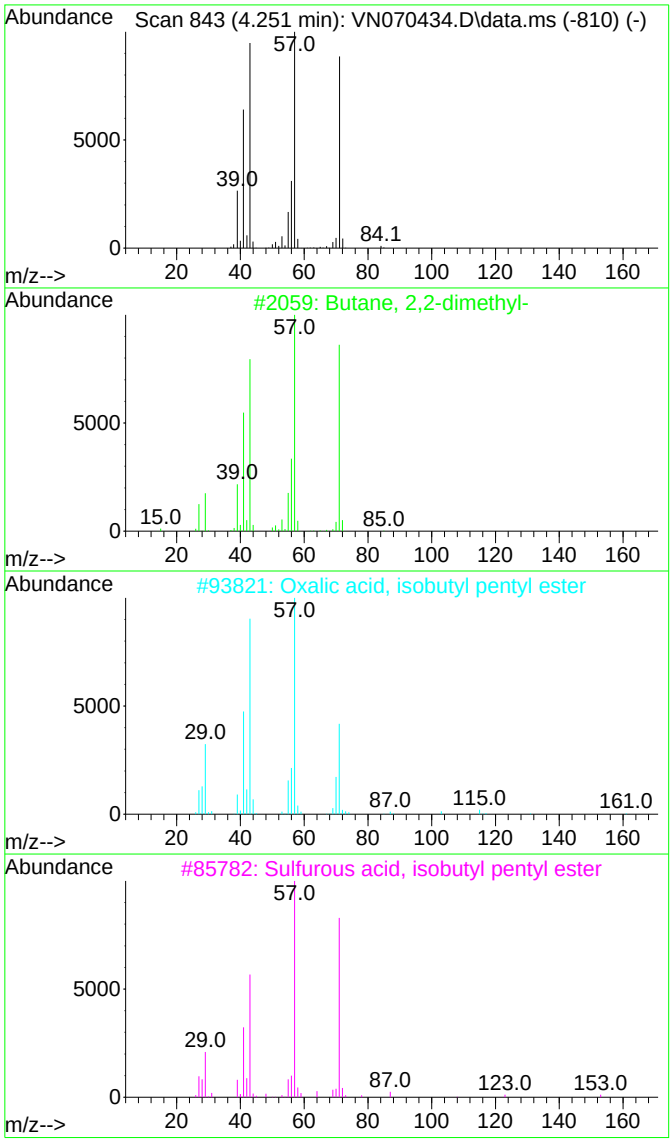
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

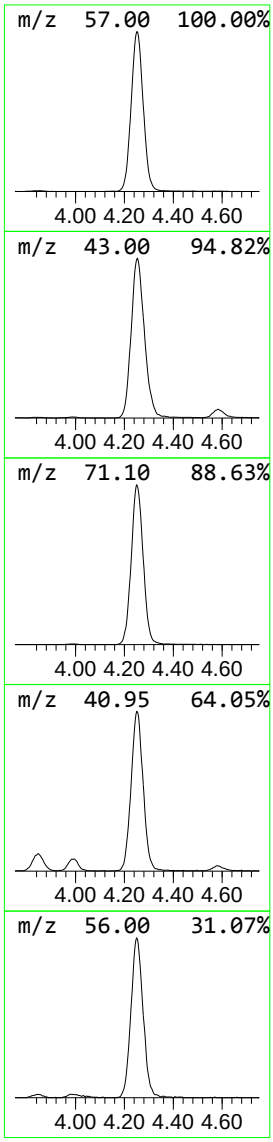
****

Peak Number 2 Butane, 2,2-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.251	38.39 ug/l	2984950	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	90
2	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	64
3	Sulfurous acid, isobutyl pentyl ...	208	C9H20O3S	1000309-13-8	56
4	Propanoic acid, 2-methyl-, anhyd...	158	C8H14O3	000097-72-3	50
5	Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

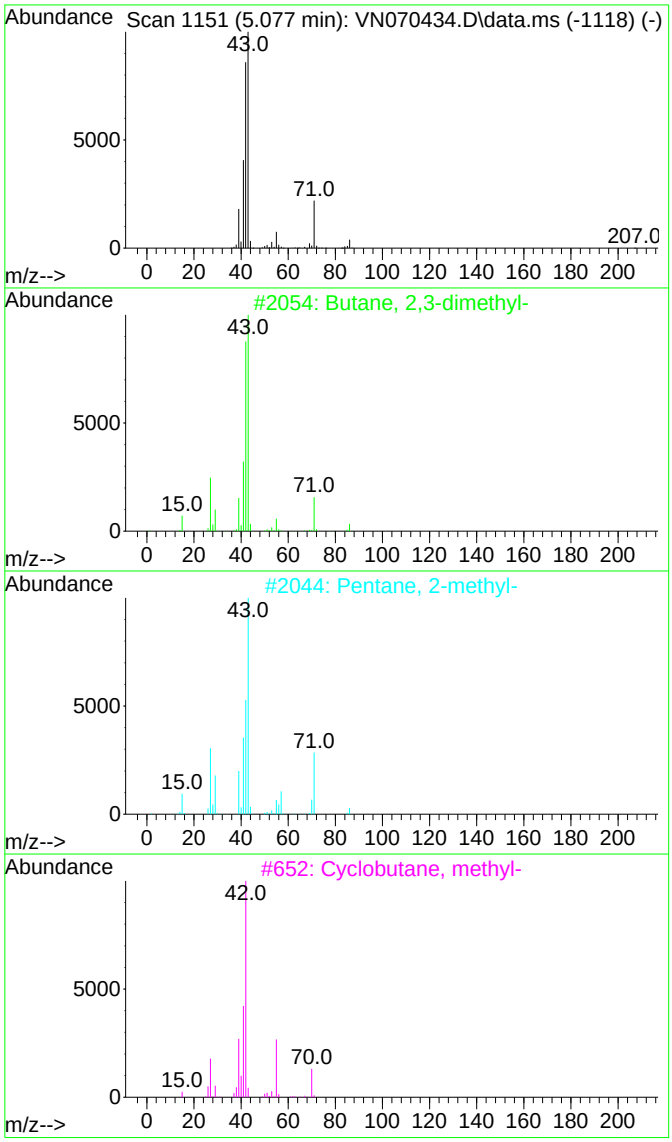
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

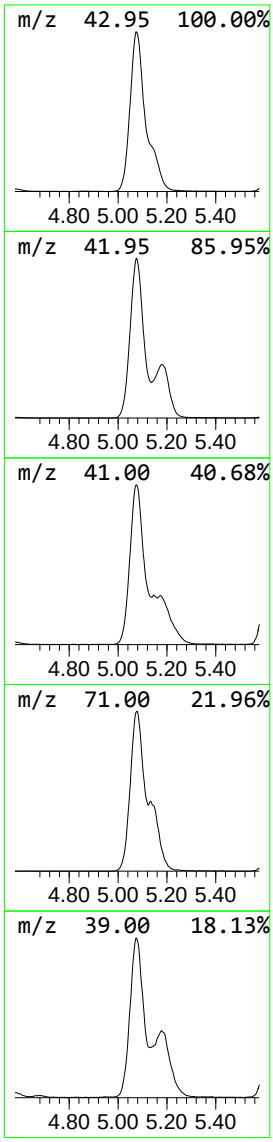
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.077	97.08 ug/l	7548190	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	91
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	50
3	Cyclobutane, methyl-	70	C5H10	000598-61-8	38
4	Pentane	72	C5H12	000109-66-0	38
5	Tetrahydrofuran	72	C4H8O	000109-99-9	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

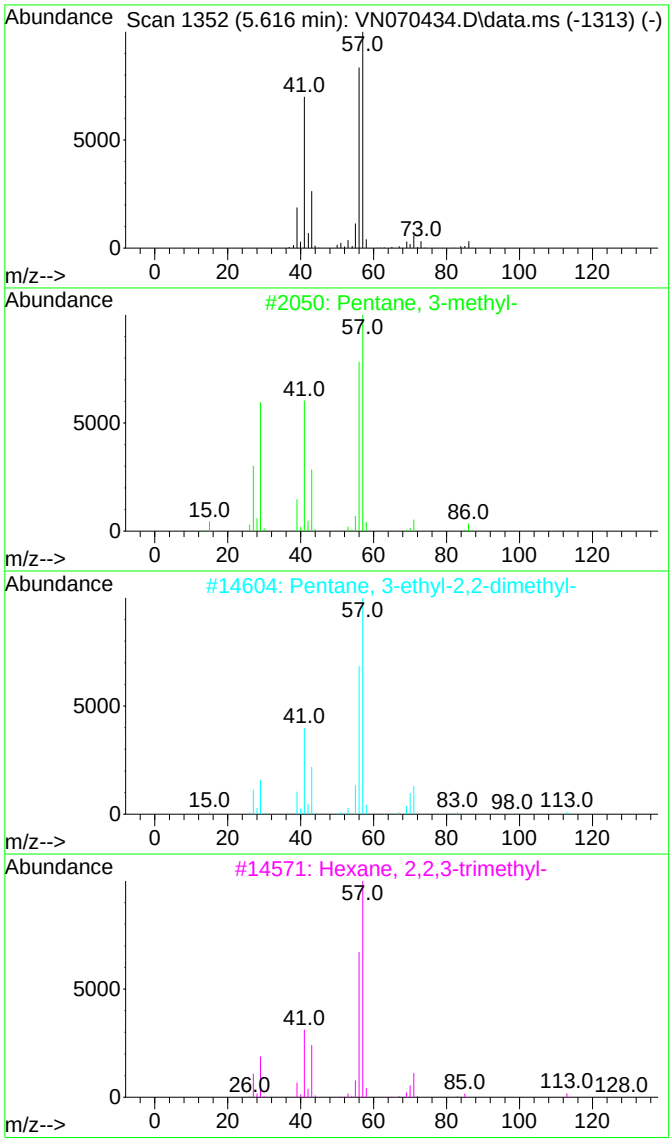
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

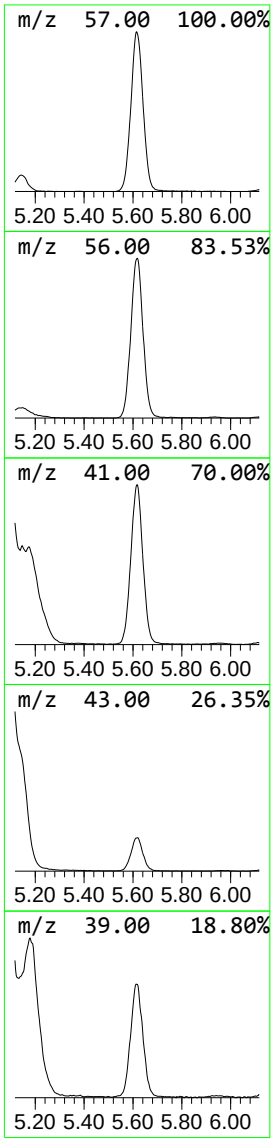
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Pentane, 3-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.616	28.60 ug/l	2223710	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	40
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	38
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	38





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

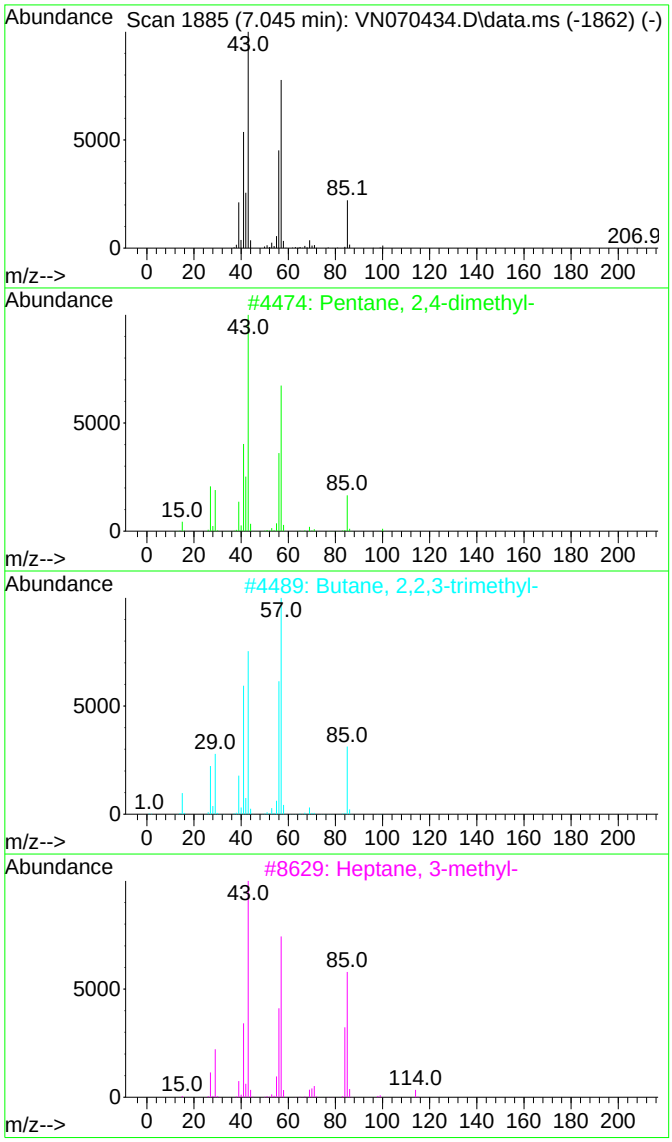
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

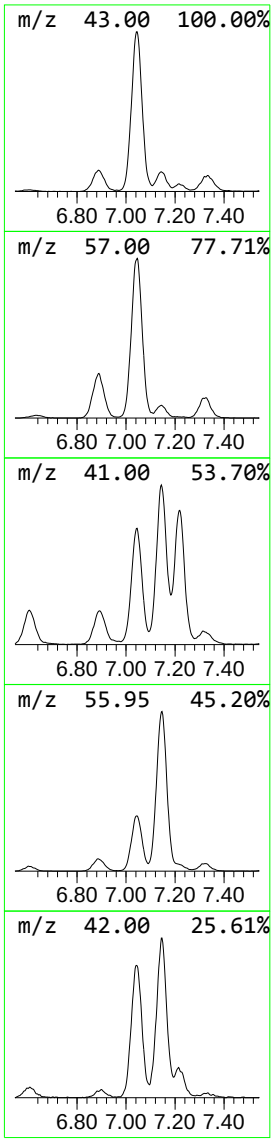
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Pentane, 2,4-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.045	23.68 ug/l	1840970	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
3	Heptane, 3-methyl-	114	C8H18	000589-81-1	64
4	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

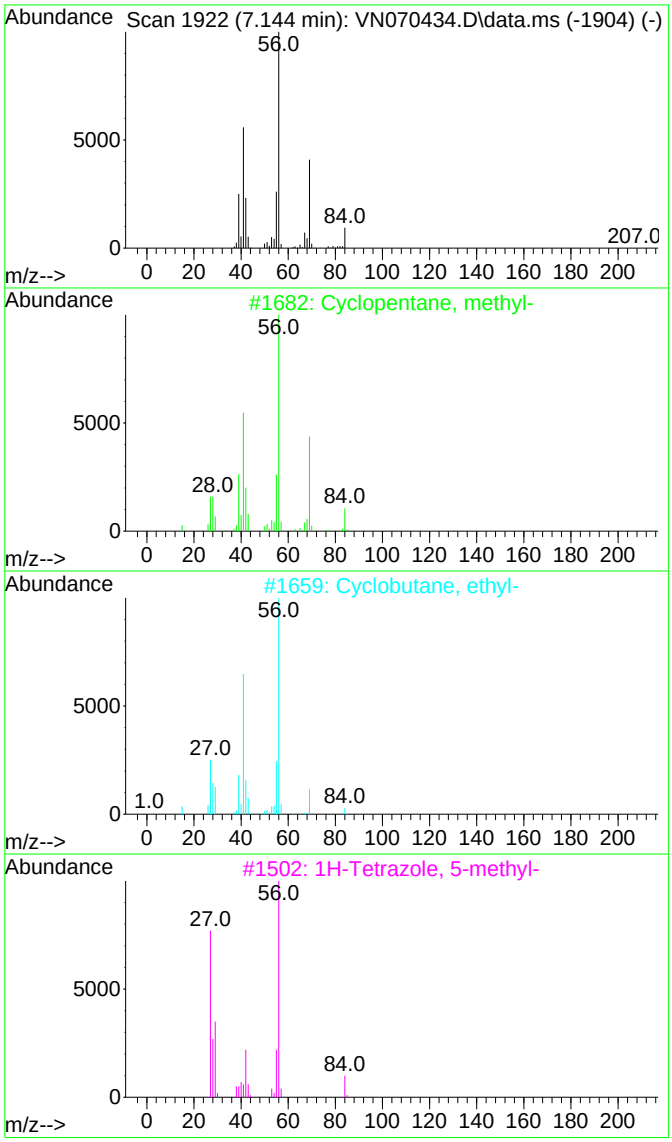
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

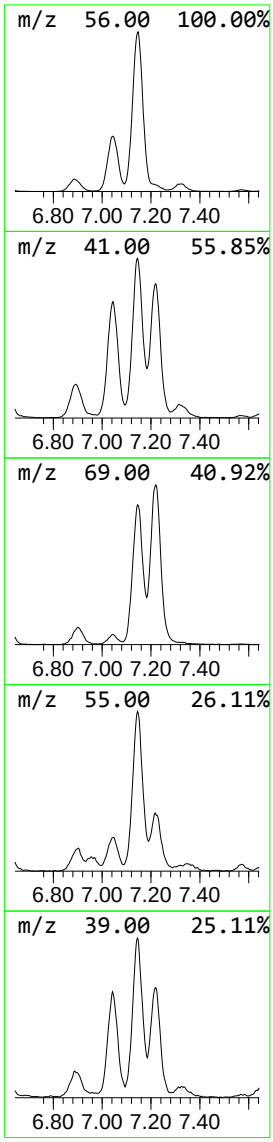
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclopentane, methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.144	27.52 ug/l	2139760	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
4	Cyclobutane	56	C4H8	000287-23-0	53
5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	39





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Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

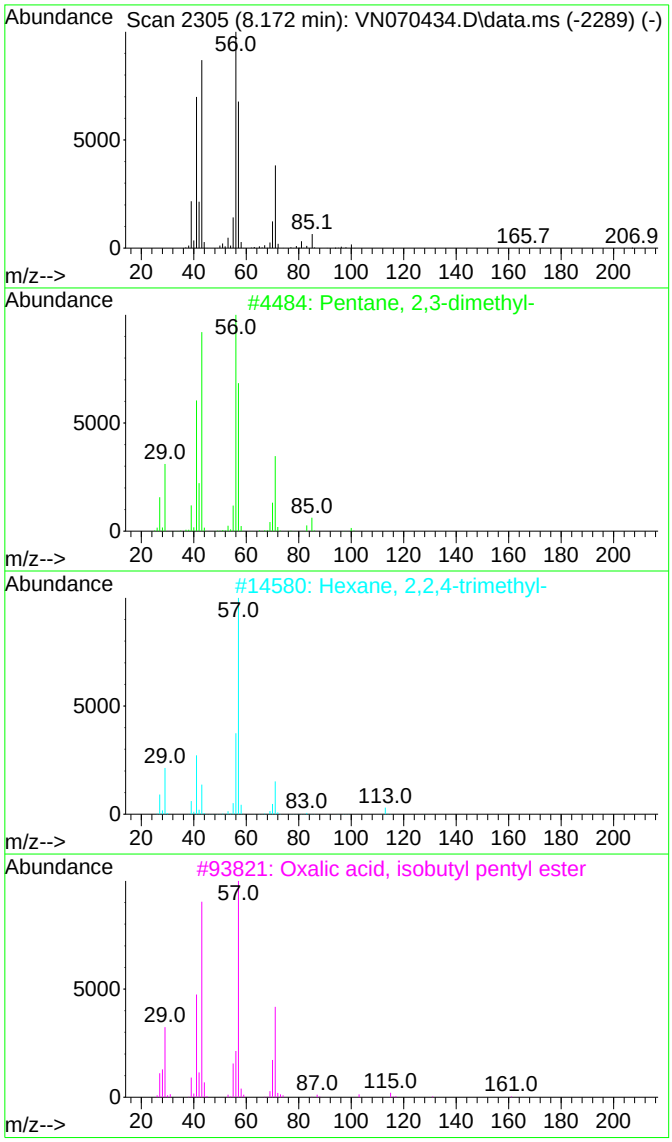
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

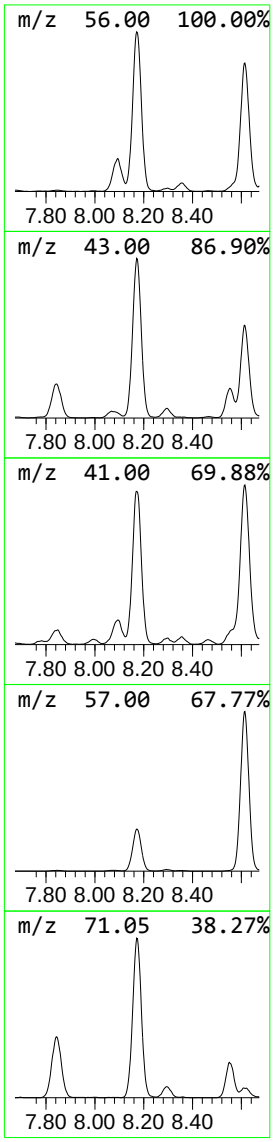
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentane, 2,3-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.172	58.48 ug/l	4547220	Pentafluorobenzene	8.083

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	42
3	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	38
4	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	38
5	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

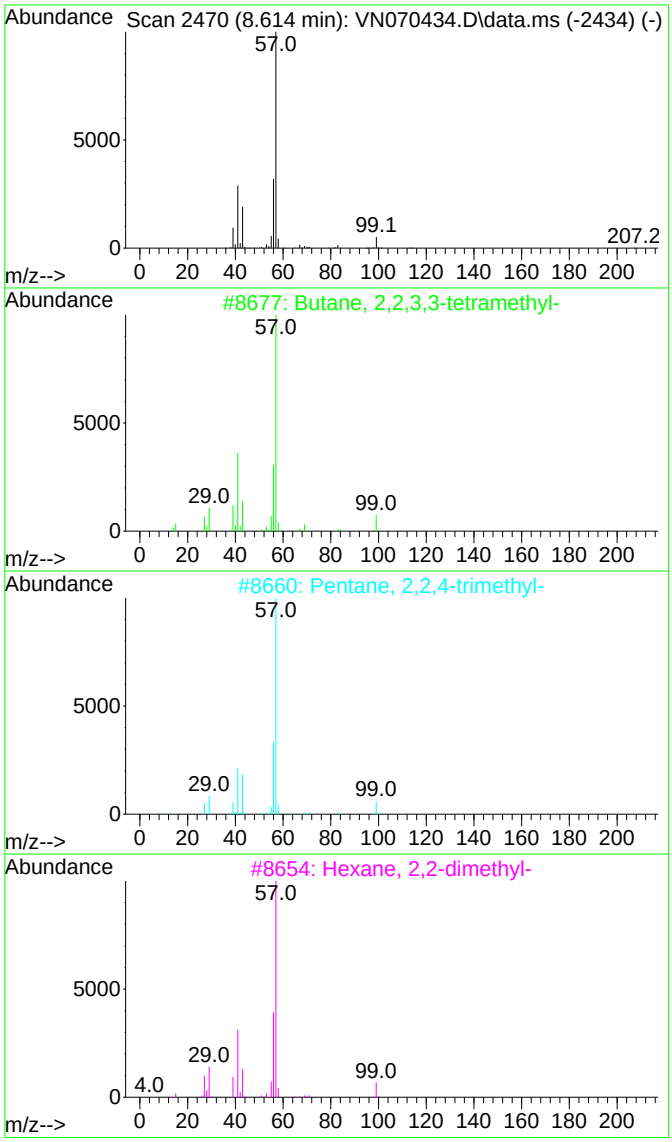
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

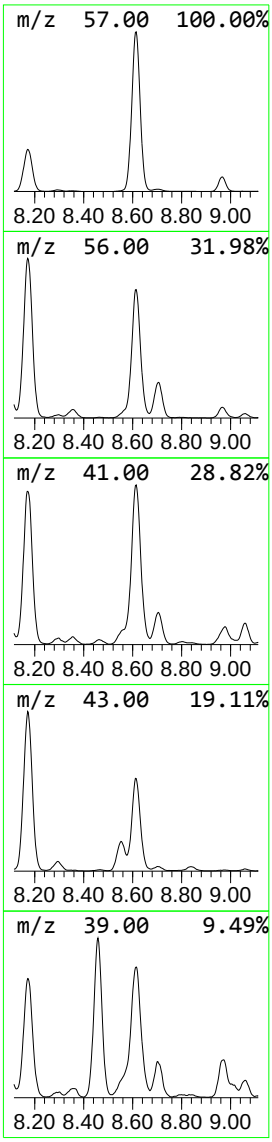
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.614	61.91 ug/l	6248810	1,4-Difluorobenzene	8.965

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	78
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	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	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

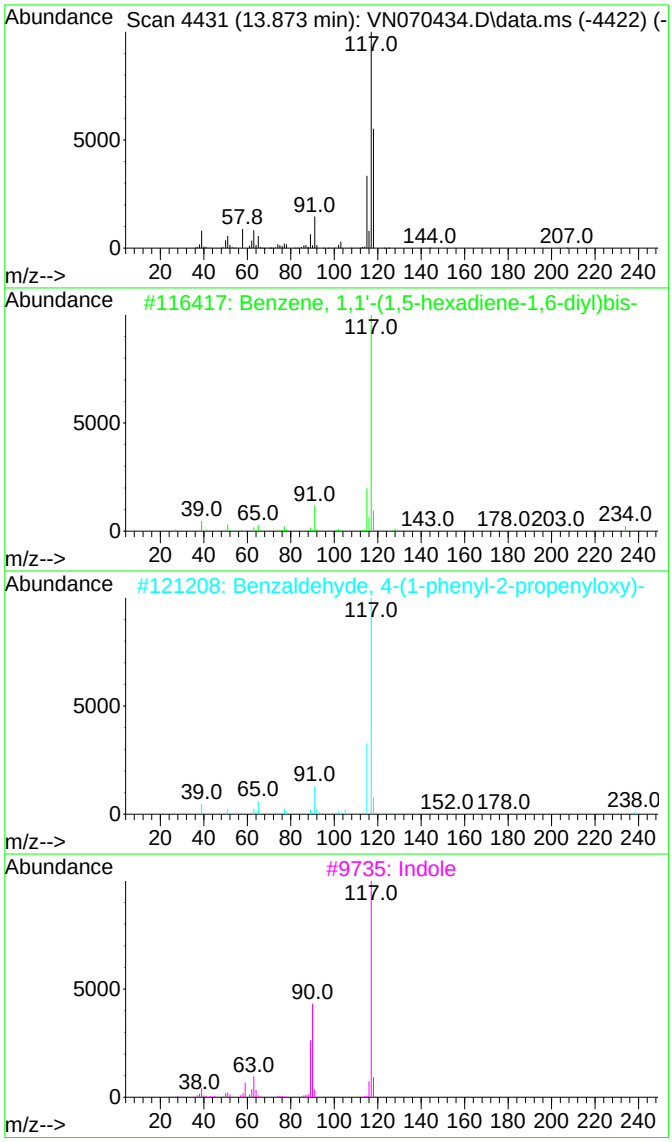
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

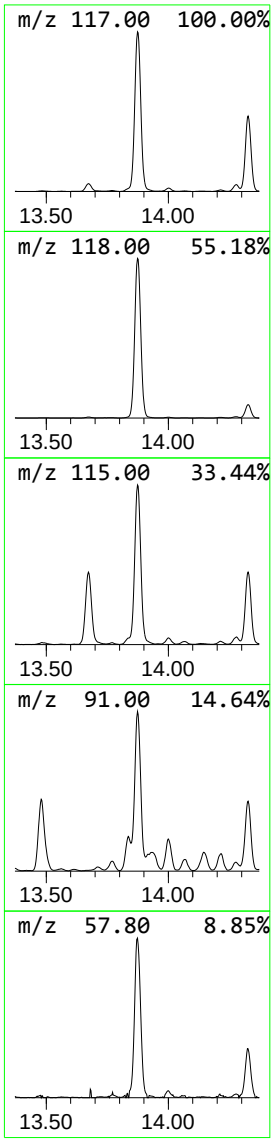
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,1'-(1,5-hexadien... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.873	84.15 ug/l	6509170	1,4-Dichlorobenzene-d4	13.672

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	Benzaldehyde, 4-(1-phenyl-2-prop...	238	C16H14O2	1000277-56-1	53
3	Indole	117	C8H7N	000120-72-9	49
4	Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	45
5	4-Ethynylaniline	117	C8H7N	014235-81-5	43





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Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

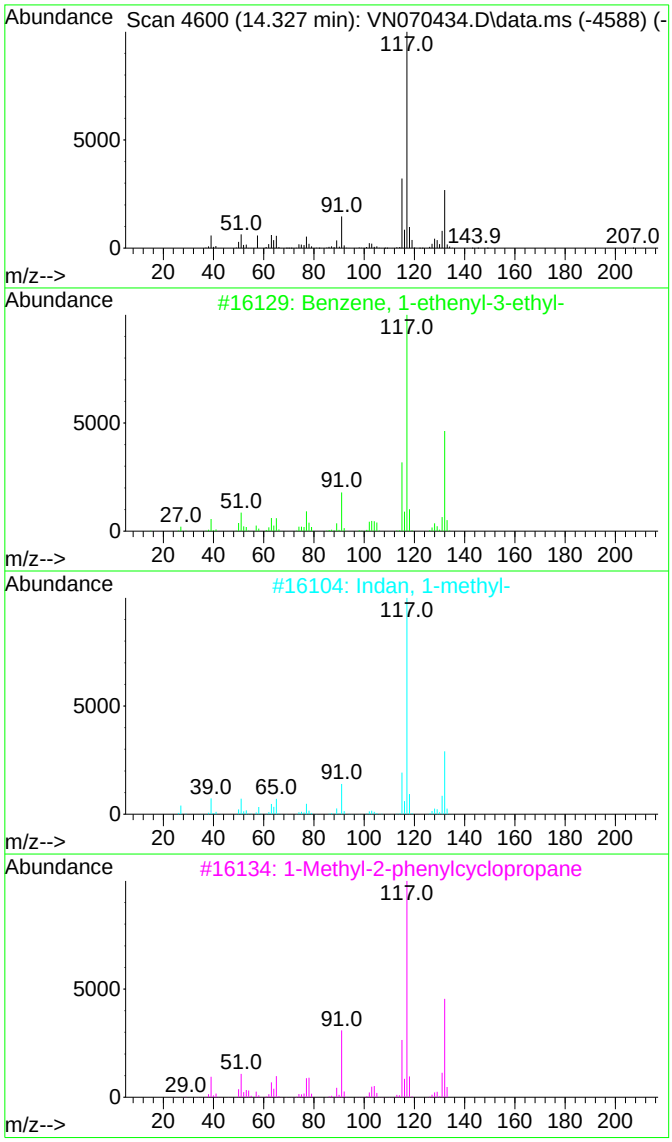
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

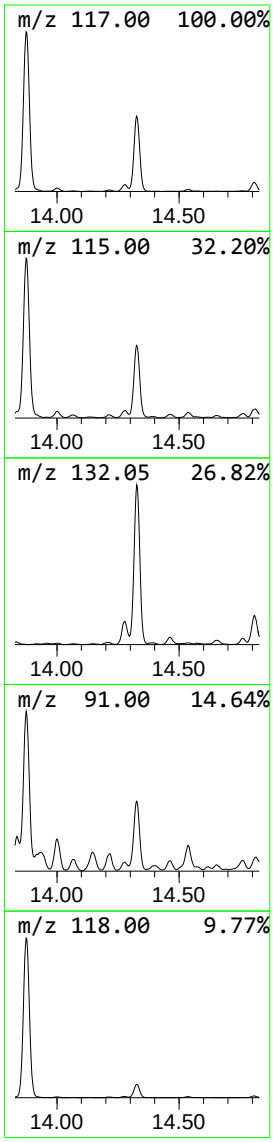
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.327	40.48 ug/l	3131090	1,4-Dichlorobenzene-d4	13.672

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	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	72
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	72





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070434.D
Acq On : 3 Jan 2022 13:54
Operator : JC/MD
Sample : M5212-06
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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
Butane, 2-methyl-	3.138	139.0	ug/l	10806400	1	8.083	3887550	50.0
Butane, 2,2-dim...	4.251	38.4	ug/l	2984950	1	8.083	3887550	50.0
Butane, 2,3-dim...	5.077	97.1	ug/l	7548190	1	8.083	3887550	50.0
Pentane, 3-methyl-	5.616	28.6	ug/l	2223710	1	8.083	3887550	50.0
Pentane, 2,4-di...	7.045	23.7	ug/l	1840970	1	8.083	3887550	50.0
Cyclopentane, m...	7.144	27.5	ug/l	2139760	1	8.083	3887550	50.0
Pentane, 2,3-di...	8.172	58.5	ug/l	4547220	1	8.083	3887550	50.0
Butane, 2,2,3,3...	8.614	61.9	ug/l	6248810	2	8.965	5046490	50.0
Benzene, 1,1'-(...	13.873	84.2	ug/l	6509170	4	13.672	3867610	50.0
Benzene, 1-ethe...	14.327	40.5	ug/l	3131090	4	13.672	3867610	50.0