LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 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: VN070433.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	113	rBV	418454	792559	7.56%	0.640%
2	2.445	159	170	188	rVB	310840	502761	4.79%	0.406%
3	3.140	408	429	462	rBV	4539871	10489421	100.00%	8.467%
4	3.516	554	569	595	rVB2	92861	232520	2.22%	0.188%
5	3.993	725	747	768	rBV3	69429	191708	1.83%	0.155%
6	4.253	811	844	881	rBV3	845112	2997891	28.58%	2.420%
7	5.076	1117	1151	1174	rBV2	1889687	7580349	72.27%	6.119%
8	5.618	1316	1353	1395	rBV	605874	2120693	20.22%	1.712%
9	6.605	1700	1721	1747	rBV5	102341	314741	3.00%	0.254%
10	6.892	1797	1828	1855	rBV5	158154	557532	5.32%	0.450%
11	7.045	1862	1885	1904	rBV3	646616	1928411	18.38%	1.557%
12	7.144	1905	1922	1937	rVV2	611627	1772978	16.90%	1.431%
13	7.219	1940	1950	1973	rVB3	363666	932680	8.89%	0.753%
14	7.324	1978	1989	2019	rVB9	70652	232116	2.21%	0.187%
15	7.651	2093	2111	2141	rVB4	56167	178637	1.70%	0.144%
16	7.785	2141	2161	2164	rBV4	57530	118098	1.13%	0.095%
17	7.844	2168	2183	2213	rVB3	308361	858428	8.18%	0.693%
18	8.019	2223	2248	2259	rBV	580246	1441313	13.74%	1.163%
19	8.086	2260	2273	2289	rVV2	1468692	3532046	33.67%	2.851%
20	8.174	2289	2306	2335	rVV	1838727	4693738	44.75%	3.789%
21	8.295	2336	2351	2361	rVV3	72020	168318	1.60%	0.136%
22	8.354	2362	2373	2387	rVV3	111057	266520	2.54%	0.215%
23	8.456	2387	2411	2433	rVV3	2600356	7091458	67.61%	5.724%
24	8.560	2435	2450	2451	rVV4	265343	454460	4.33%	0.367%
25	8.614	2455	2470	2491	rVV	2190576	5733073	54.66%	4.628%
26	8.705	2492	2504	2527	rVB2	433631	988901	9.43%	0.798%
27	8.965	2579	2601	2626	rBV	2023844	4585567	43.72%	3.702%
28	9.059	2627	2636	2652	rVB2	194123	393445	3.75%	0.318%
29	9.199	2672	2688	2697	rBV2	260542	546980	5.21%	0.442%
30	9.244	2698	2705	2725	rVB	205504	413002	3.94%	0.333%
31	9.421	2750	2771	2795	rBV4	658763	2079158	19.82%	1.678%
32	9.510	2796	2804	2822	rVB4	223404	447694	4.27%	0.361%
33	9.590	2822	2834	2844	rBV3	91740	180451	1.72%	0.146%
34	9.730	2866	2886	2901	rBV3	191506	461776	4.40%	0.373%
35	9.880	2929	2942	2962	rVB	959350	1964346	18.73%	1.586%
36	9.974	2963	2977	2979	rVV2	269834	383908	3.66%	0.310%

LSC Area Percent Report

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Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 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	10.014	2981	2992	3019	rVV	999369	2284157	21.78%	1.844%
38	10.121	3021	3032	3045	rVV7	88451	217094	2.07%	0.175%
39	10.196	3046	3060	3074	rVV5	198648	421158	4.02%	0.340%
40	10.263	3075	3085	3093	rVV	68054	126512	1.21%	0.102%
41	10.320	3094	3106	3116	rVV2	387066	773142	7.37%	0.624%
42	10.365	3117	3123	3135	rVV3	236365	409270	3.90%	0.330%
43	10.440	3135	3151	3195	rVV	2629585	5454203	52.00%	4.403%
44	10.639	3198	3225	3240	rVV3	1034118	2365431	22.55%	1.909%
45	10.714	3240	3253	3267	rVV3	923906	1722610	16.42%	1.391%
46	10.781	3268	3278	3289	rVV2	123919	254526	2.43%	0.205%
47	10.856	3291	3306	3328	rVV4	726705	1743670	16.62%	1.408%
48	10.942	3328	3338	3358	rVB7	85085	218213	2.08%	0.176%
49	11.127	3400	3407	3417	rVB3	113078	170201	1.62%	0.137%
50	11.256	3447	3455	3466	rVB5	147431	244772	2.33%	0.198%
51	11.449	3518	3527	3543	rVB3	131021	226489	2.16%	0.183%
52	11.556	3558	3567	3585	rVB10	46855	116896	1.11%	0.094%
53	11.744	3621	3637	3652	rBV	2503902	4567846	43.55%	3.687%
54	11.843	3661	3674	3698	rVB	2870101	5149367	49.09%	4.157%
55	12.578	3934	3948	3969	rBV	1690816	2824580	26.93%	2.280%
56	12.731	3990	4005	4026	rVB	1819715	3213468	30.64%	2.594%
57	12.919	4061	4075	4094	rVB	1055557	1820037	17.35%	1.469%
58	13.219	4174	4187	4203	rBV	266490	459182	4.38%	0.371%
59	13.321	4210	4225	4232	rBV3	57159	112219	1.07%	0.091%
60	13.366	4233	4242	4256	rVB	110608	188781	1.80%	0.152%
61	13.484	4273	4286	4305	rVB3	304892	710708	6.78%	0.574%
62	13.672	4342	4356	4368	rBV	1921221	3482943	33.20%	2.812%
63	13.769	4384	4392	4404	rVB2	132063	209996	2.00%	0.170%
64	13.836	4404	4417	4421	rBV	738964	1044048	9.95%	0.843%
65	13.876	4422	4432	4444	rVB	3430702	5668695	54.04%	4.576%
66	14.002	4466	4479	4491	rVB	613647	1015721	9.68%	0.820%
67	14.066	4491	4503	4516	rVB	436029	707514	6.75%	0.571%
68	14.147	4518	4533	4546	rVB2	374703	689768	6.58%	0.557%
69	14.214	4547	4558	4570	rBV2	296816	478400	4.56%	0.386%
70	14.278	4571	4582	4587	rBV	180695	271737	2.59%	0.219%
71	14.327	4588	4600	4614	rVB	1737219	2849190	27.16%	2.300%
72	14.404	4617	4629	4641	rVB5	188777	347565	3.31%	0.281%
73	14.463	4641	4651	4661	rBV	161201	238896	2.28%	0.193%
74	14.536	4662	4678	4689	rBV	478008	817572	7.79%	0.660%
75	14.654	4715	4722	4733	rVB2	101231	145012	1.38%	0.117%
76	14.761	4751	4762	4771	rBV3	163294	254204	2.42%	0.205%

LSC Area Percent Report

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Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 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	14.812	4771	4781	4795	rVB6	262279	475450	4.53%	0.384%
78	14.927	4807	4824	4837	rBV3	261348	603543	5.75%	0.487%
79	14.997	4838	4850	4862	rVB4	132032	230917	2.20%	0.186%
80	15.080	4872	4881	4890	rBV3	90254	147761	1.41%	0.119%
81	15.193	4910	4923	4934	rBV4	181189	344716	3.29%	0.278%
82	15.504	5031	5039	5050	rVB2	71511	112336	1.07%	0.091%
83	15.802	5135	5150	5163	rBV2	58556	126242	1.20%	0.102%
84	16.107	5251	5264	5279	rBV2	92069	197191	1.88%	0.159%

Sum of corrected areas: 123881626

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
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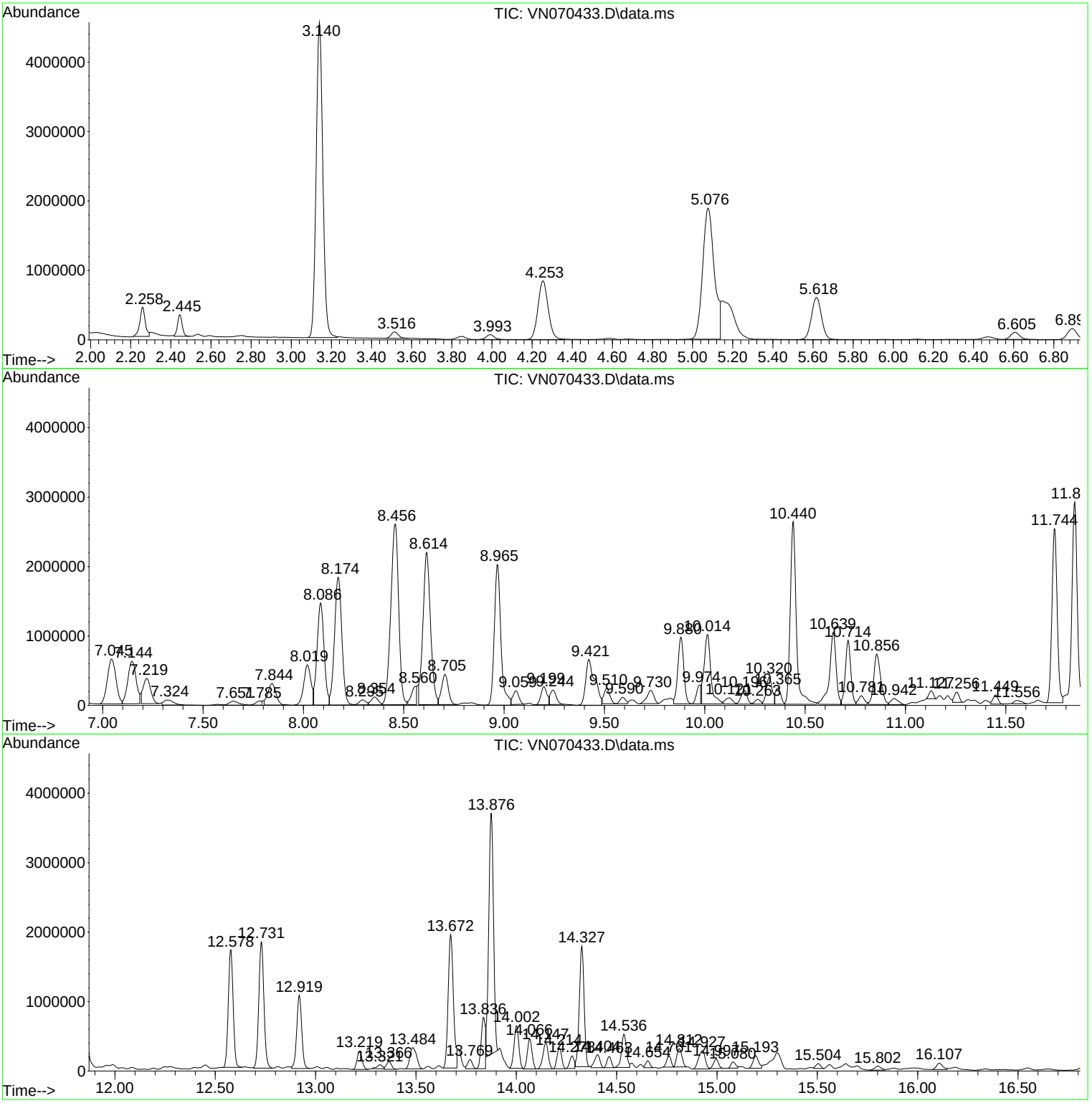
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

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 : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

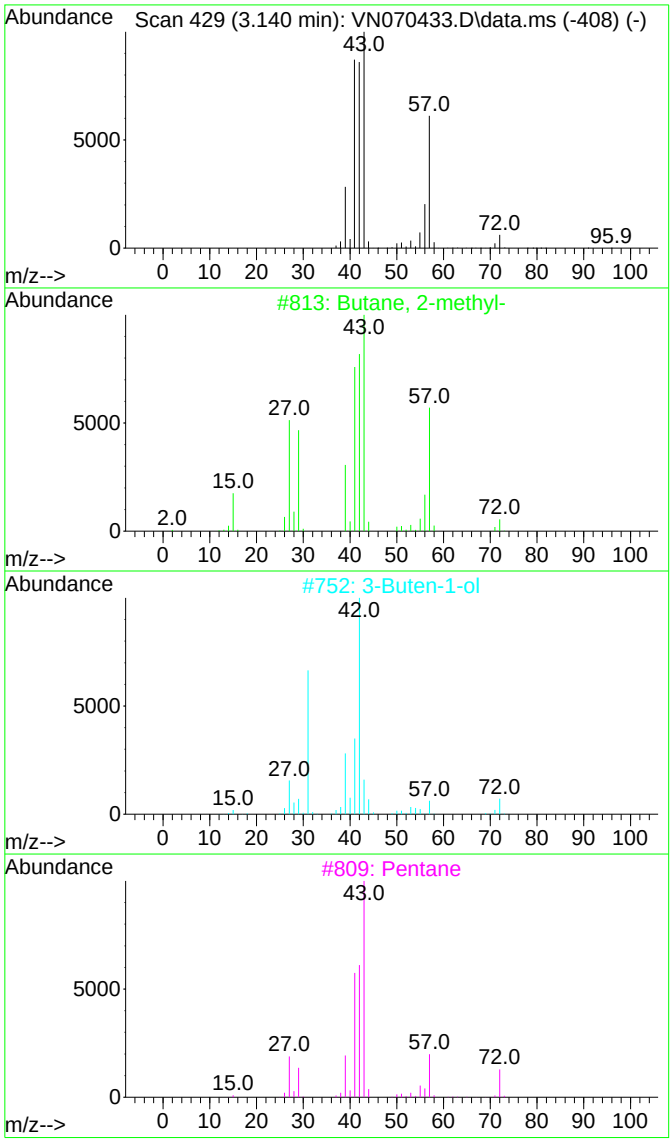
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

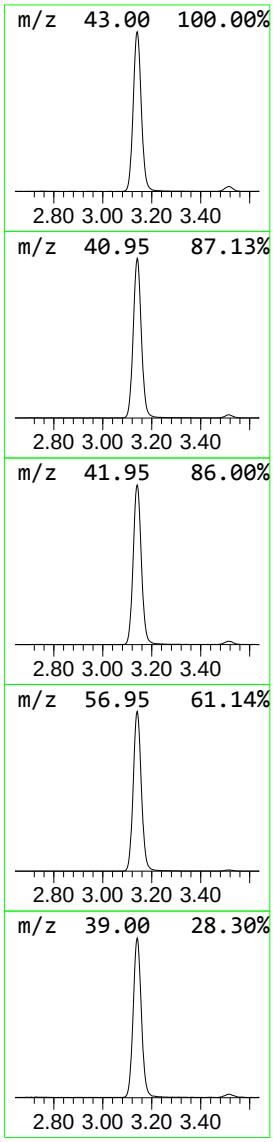
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.140	148.49 ug/l	10489400	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2	3-Buten-1-ol	72	C4H8O	000627-27-0	47
3	Pentane	72	C5H12	000109-66-0	45
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5	1-Butene	56	C4H8	000106-98-9	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

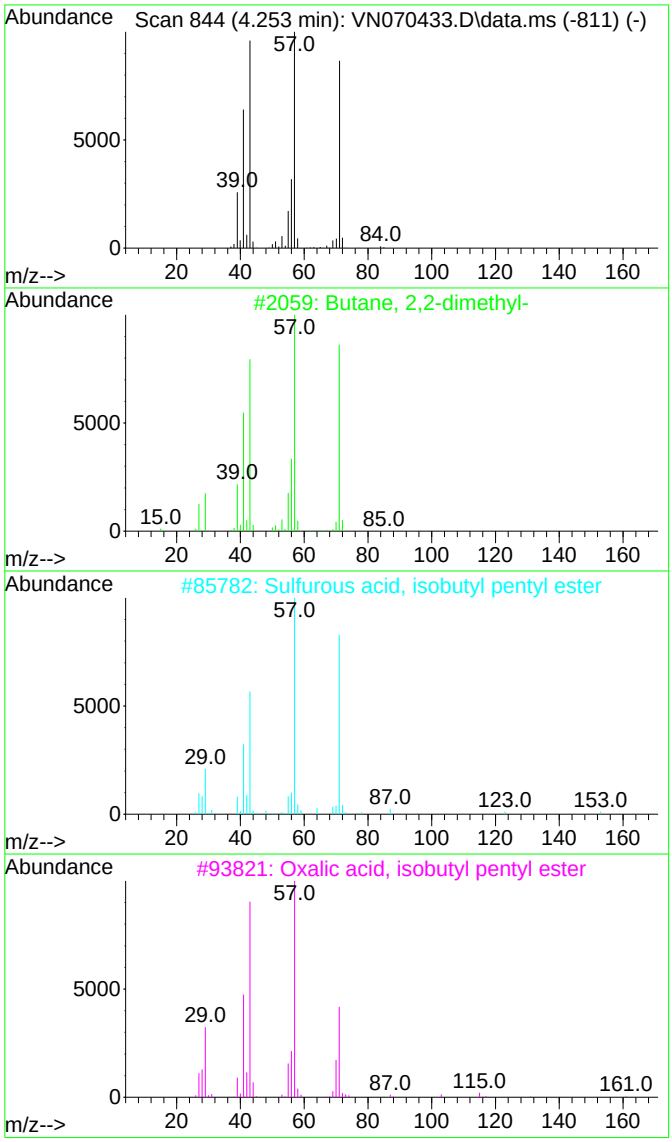
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

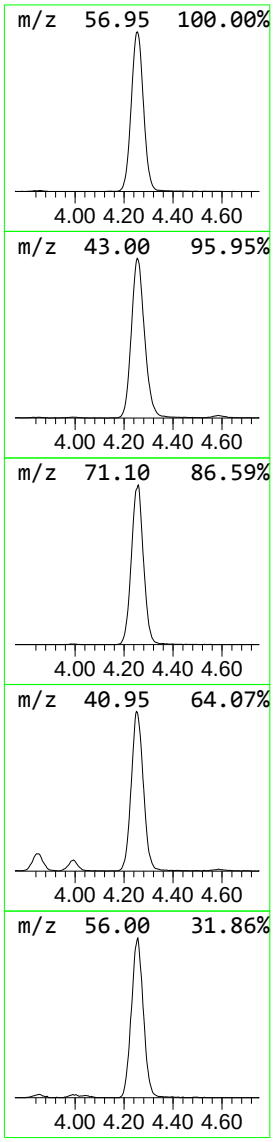
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.253	42.44 ug/l	2997890	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	90
2	Sulfurous acid, isobutyl pentyl ...	208	C9H20O3S	1000309-13-8	64
3	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	53
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	45
5	Butane, 2-iodo-3-methyl-	198	C5H11I	018295-27-7	42





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

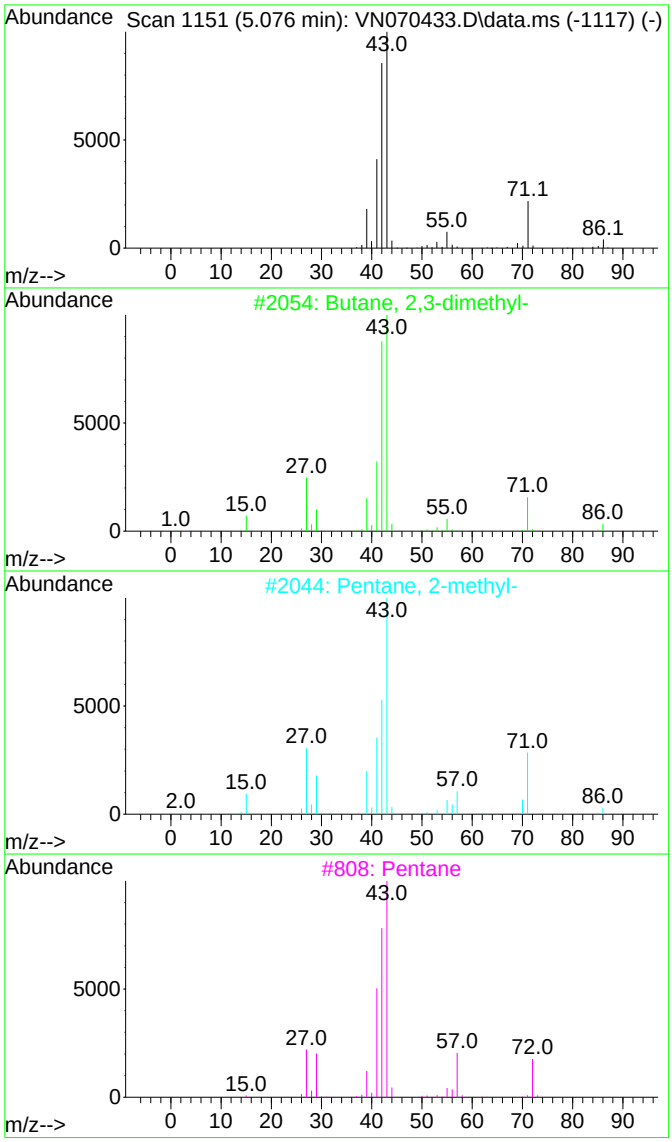
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

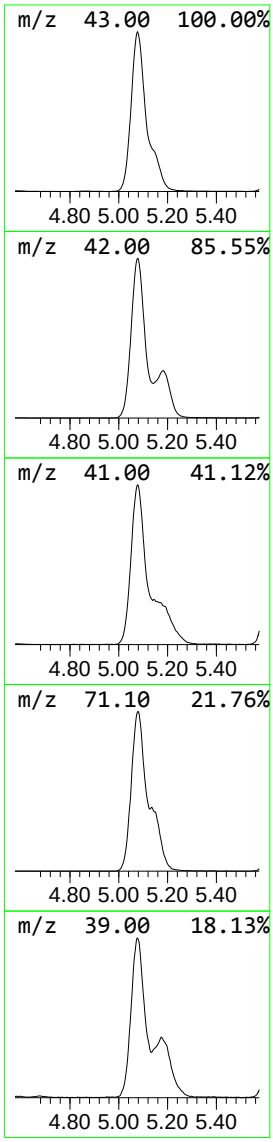
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.076	107.31 ug/l	7580350	Pentafluorobenzene	8.086

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	Pentane	72	C5H12	000109-66-0	38
4	Pyrrolidine	71	C4H9N	000123-75-1	9
5	Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	9





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ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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

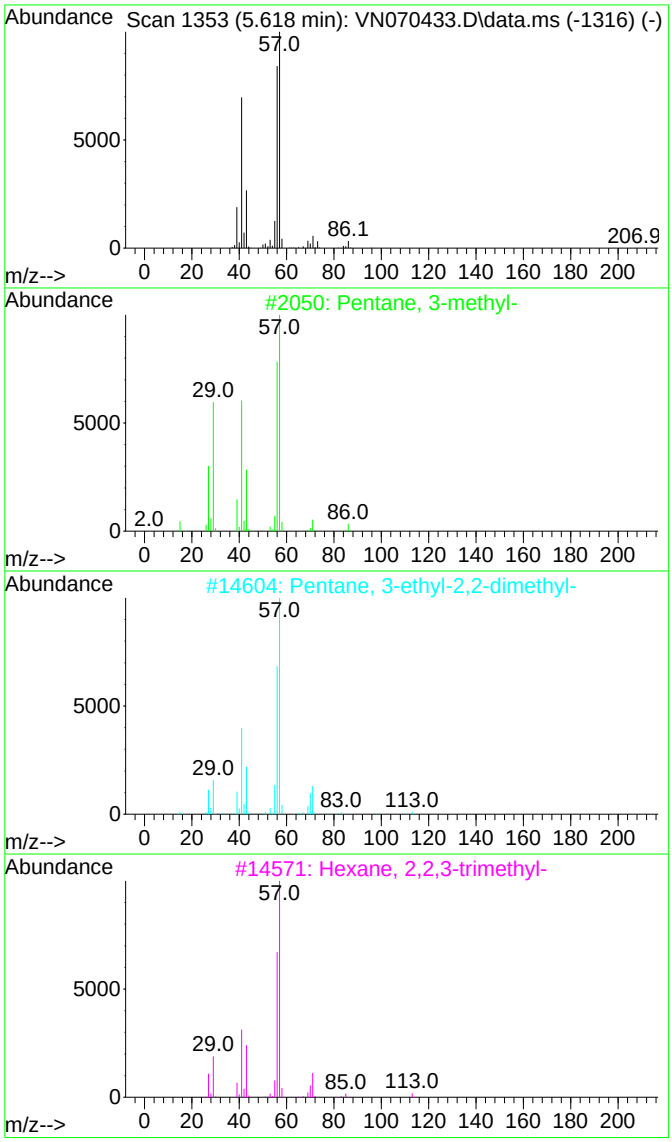
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

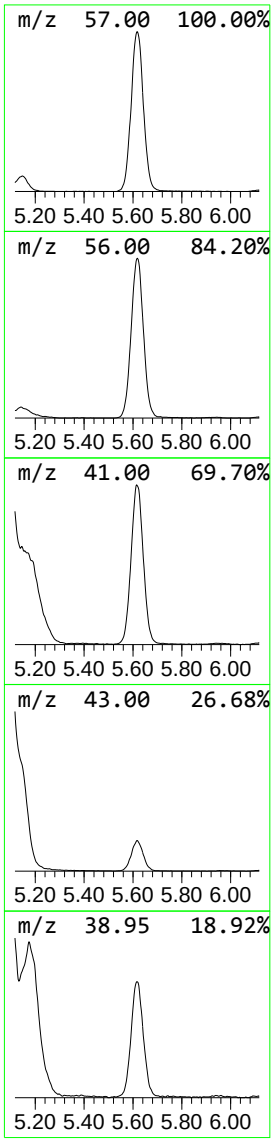
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.618	30.02 ug/l	2120690	Pentafluorobenzene	8.086

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	78
3	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56
4	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

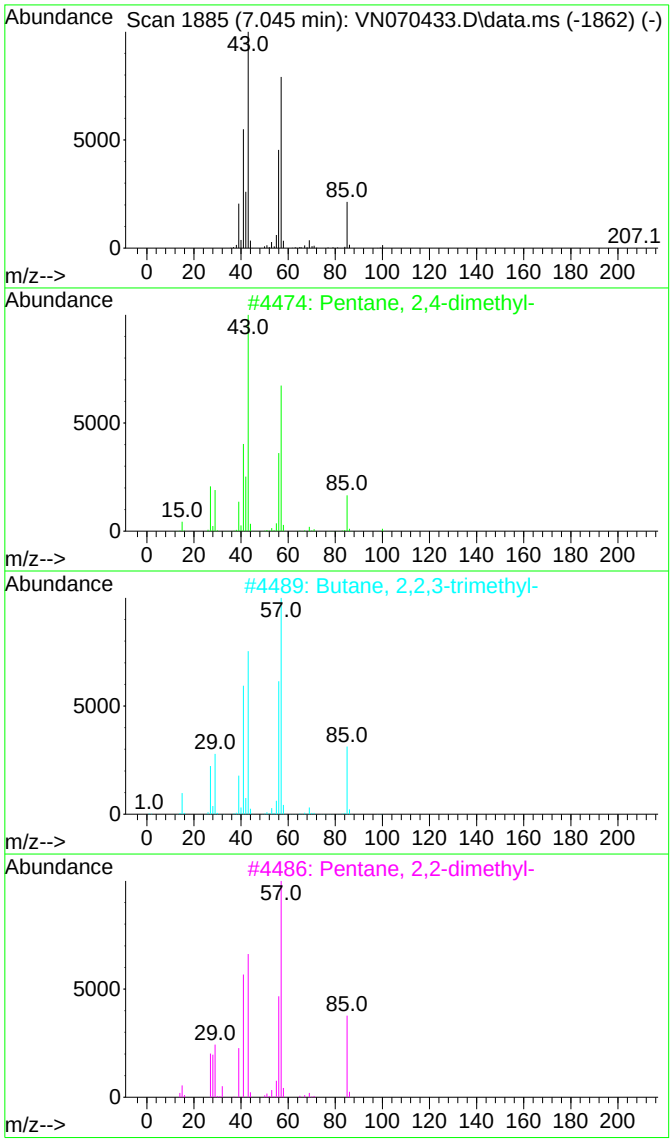
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

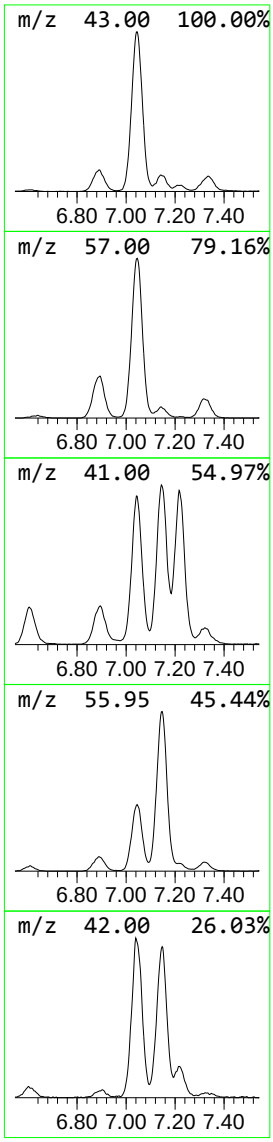
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Pentane, 2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.045	27.30 ug/l	1928410	Pentafluorobenzene	8.086

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	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	50
4	2,2,6,6-Tetramethylheptane	156	C11H24	040117-45-1	50
5	Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	45





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

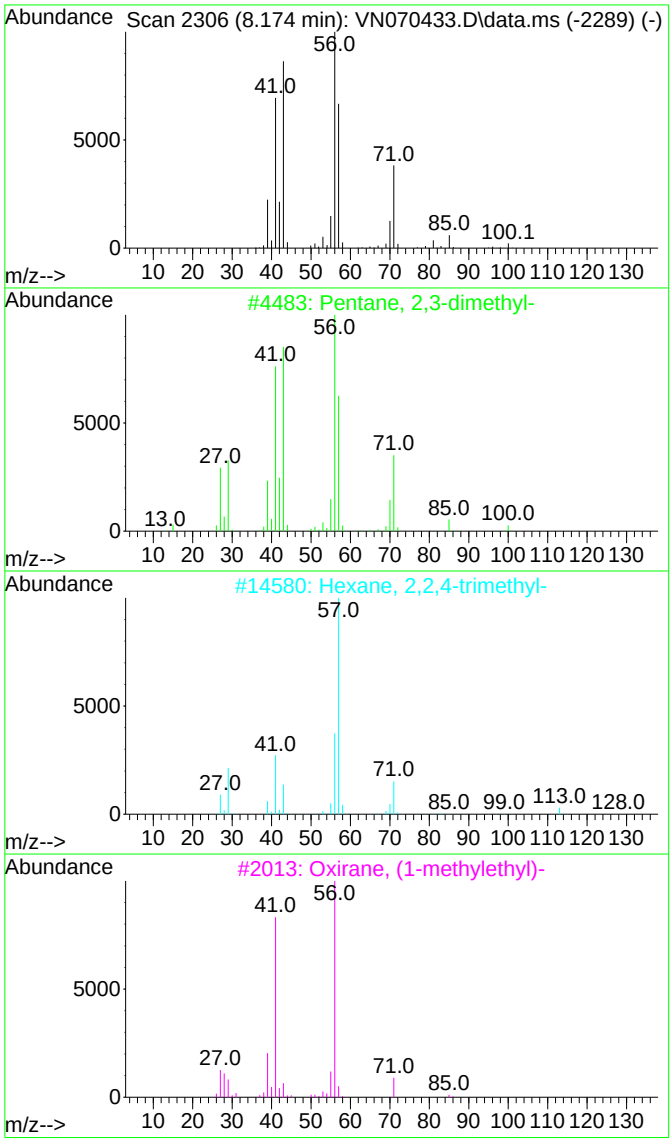
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

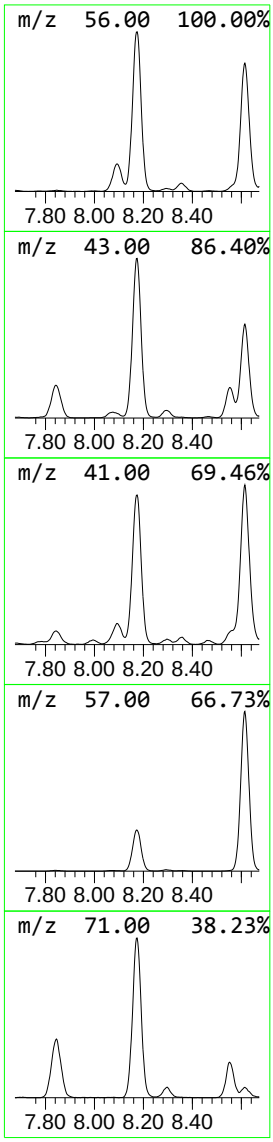
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Pentane, 2,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.174	66.44 ug/l	4693740	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	95
2	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	42
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	38
4	Oxalic acid, isobutyl pentyl ester	216	C11H20O4	1000309-37-0	38
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

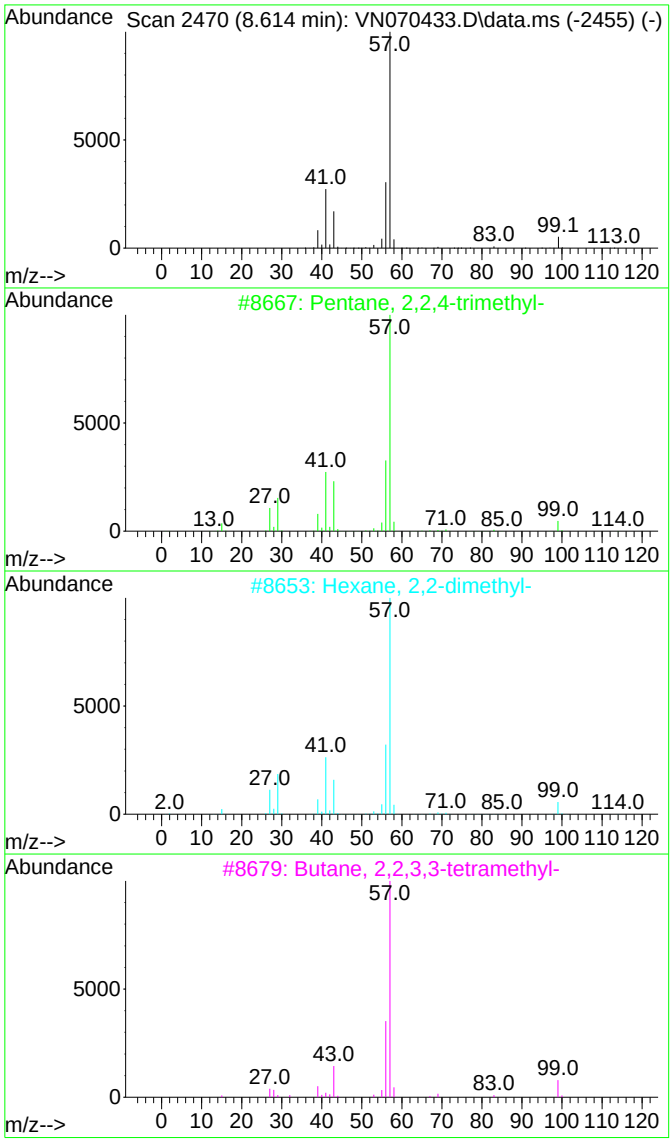
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

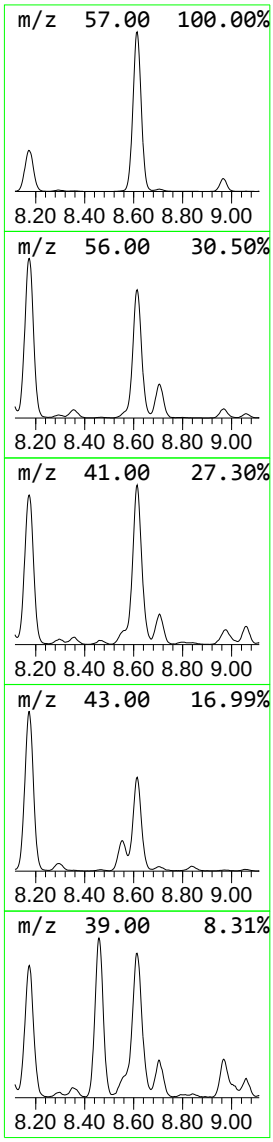
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Pentane, 2,2,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.614	62.51 ug/l	5733070	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	83
2	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	83
3	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	78
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	64





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Data File : VN070433.D
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Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

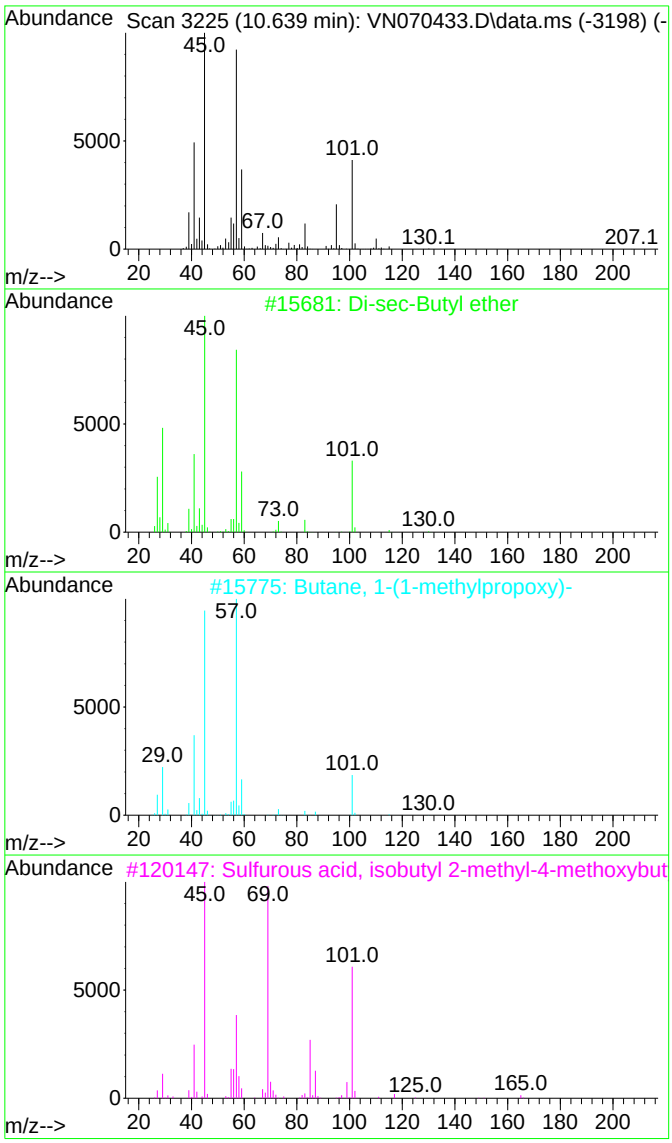
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

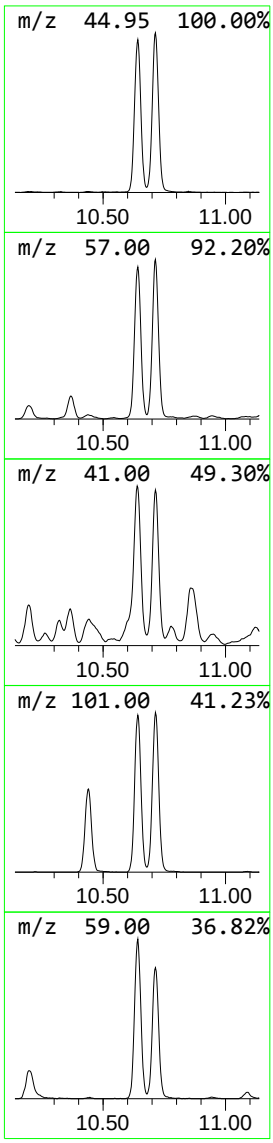
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Di-sec-Butyl ether Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.639	25.89 ug/l	2365430	Chlorobenzene-d5	11.744

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Di-sec-Butyl ether	130	C8H18O	006863-58-7	87
2	Butane, 1-(1-methylpropoxy)-	130	C8H18O	000999-65-5	50
3	Sulfurous acid, isobutyl 2-methy...	238	C10H22O4S	1000309-20-3	50
4	Boronic acid, ethyl-, diethyl ester	130	C6H15BO2	053907-92-9	38
5	1-Octene, 3-(methoxymethoxy)-	172	C10H20O2	088738-34-5	37





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Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

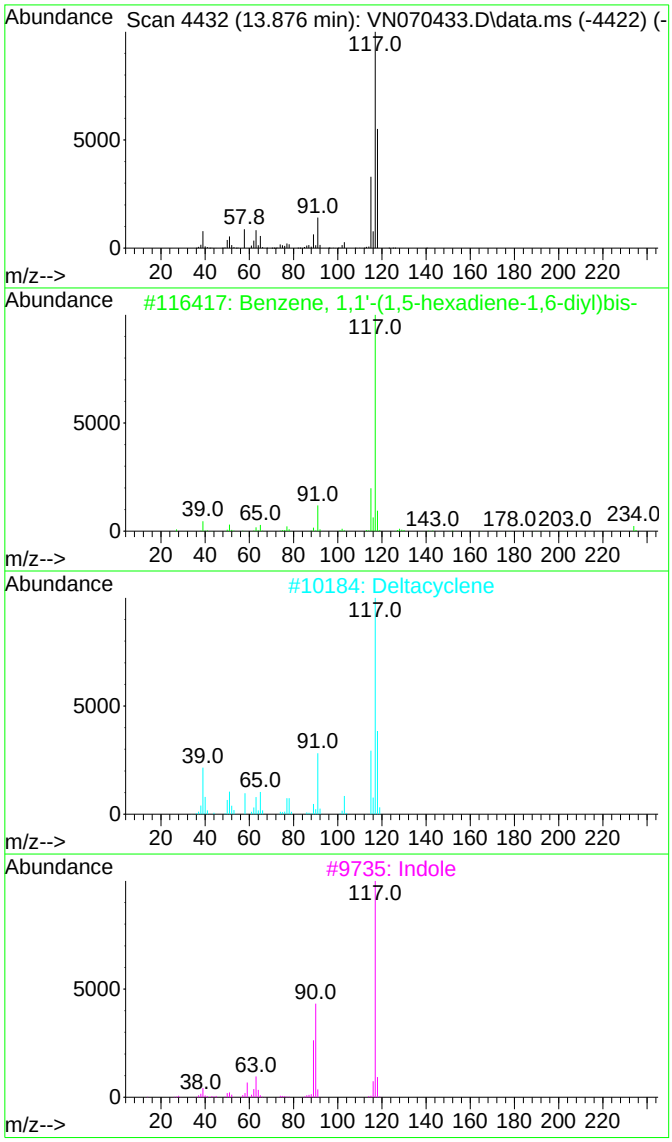
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

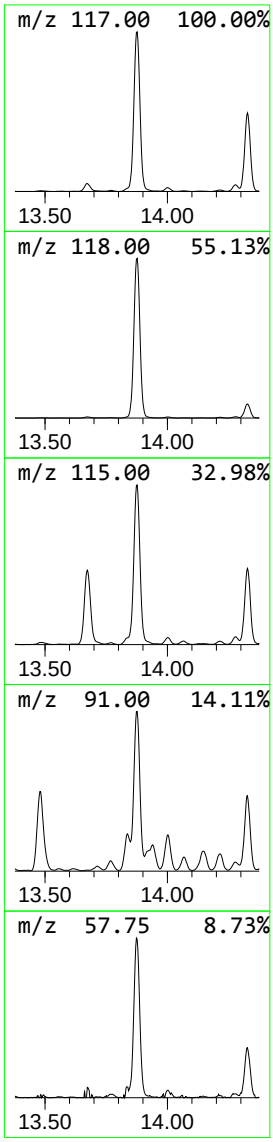
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.876	81.38 ug/l	5668700	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	Deltacyclene	118	C9H10	007785-10-6	50
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 : VN070433.D
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Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :

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

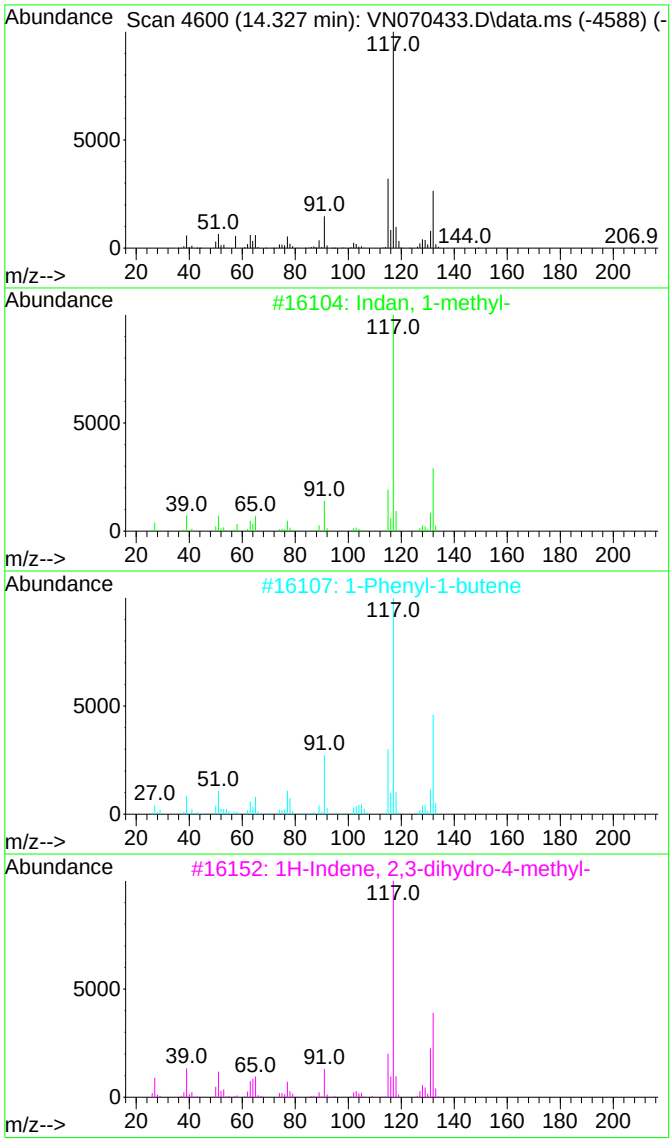
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

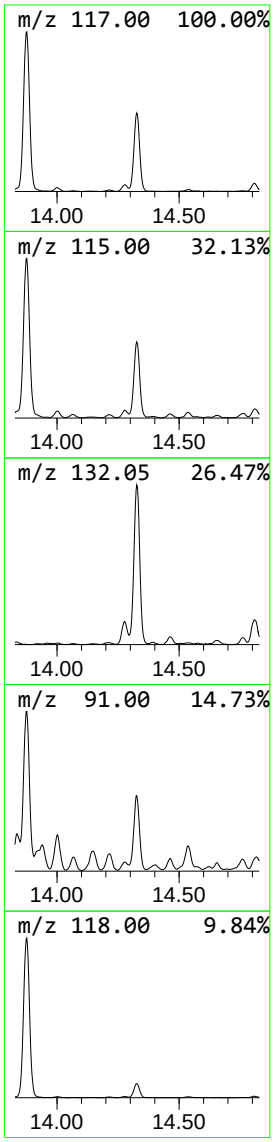
****

Peak Number 10 Indan, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.327	40.90 ug/l	2849190	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	90
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
3	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	83
4	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	80
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	78





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070433.D
Acq On : 3 Jan 2022 13:29
Operator : JC/MD
Sample : M5212-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 8 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-	3.140	148.5	ug/l	10489400	1	8.086	3532050	50.0
Butane, 2,2-dim...	4.253	42.4	ug/l	2997890	1	8.086	3532050	50.0
Butane, 2,3-dim...	5.076	107.3	ug/l	7580350	1	8.086	3532050	50.0
Pentane, 3-methyl-	5.618	30.0	ug/l	2120690	1	8.086	3532050	50.0
Pentane, 2,4-di...	7.045	27.3	ug/l	1928410	1	8.086	3532050	50.0
Pentane, 2,3-di...	8.174	66.4	ug/l	4693740	1	8.086	3532050	50.0
Pentane, 2,2,4-...	8.614	62.5	ug/l	5733070	2	8.965	4585570	50.0
Di-sec-Butyl ether	10.639	25.9	ug/l	2365430	3	11.744	4567850	50.0
Benzene, 1,1'-(...	13.876	81.4	ug/l	5668700	4	13.672	3482940	50.0
Indan, 1-methyl-	14.327	40.9	ug/l	2849190	4	13.672	3482940	50.0