LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 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\82N011822W.M
Title : SW846 8260

Signal : TIC: VN070811.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.448	152	171	192	rVB	550202	934998	3.12%	0.412%
2	3.148	411	432	466	rBV	2355641	5333419	17.82%	2.350%
3	3.521	549	571	617	rBV	3206376	7711731	25.77%	3.398%
4	3.722	627	646	674	rVB2	417838	1026367	3.43%	0.452%
5	3.998	729	749	784	rVB2	301446	825739	2.76%	0.364%
6	4.259	821	846	882	rBV5	102975	385425	1.29%	0.170%
7	4.950	1070	1104	1129	rBV5	196084	914310	3.06%	0.403%
8	5.157	1131	1181	1218	rBV6	1367233	7764337	25.94%	3.421%
9	5.621	1320	1354	1395	rBV3	995856	3435873	11.48%	1.514%
10	5.937	1445	1472	1502	rVB5	124583	415665	1.39%	0.183%
11	6.123	1508	1541	1579	rBV2	1523680	4684049	15.65%	2.064%
12	6.407	1624	1647	1656	rBV4	212467	603991	2.02%	0.266%
13	6.903	1812	1832	1862	rVB6	181937	565840	1.89%	0.249%
14	7.150	1899	1924	1969	rVB	3655799	10799642	36.09%	4.758%
15	7.844	2167	2183	2204	rVB	655952	1622784	5.42%	0.715%
16	8.021	2233	2249	2256	rBV	500177	1085324	3.63%	0.478%
17	8.094	2258	2276	2296	rVB3	5042971	13339657	44.57%	5.877%
18	8.295	2333	2351	2363	rBV	716217	1657647	5.54%	0.730%
19	8.357	2365	2374	2389	rVB2	450245	963437	3.22%	0.424%
20	8.459	2389	2412	2436	rBV	12455967	29927499	100.00%	13.185%
21	8.574	2438	2455	2469	rVV7	1297380	3254960	10.88%	1.434%
22	8.641	2470	2480	2492	rVV	954371	1947293	6.51%	0.858%
23	8.708	2492	2505	2527	rVB	1531862	3548658	11.86%	1.563%
24	8.826	2531	2549	2579	rVB	958598	2192157	7.32%	0.966%
25	8.968	2581	2602	2616	rBV2	2162315	4823541	16.12%	2.125%
26	9.242	2676	2704	2738	rVB5	191684	596112	1.99%	0.263%
27	9.459	2748	2785	2825	rBV	6399160	15959147	53.33%	7.031%
28	9.639	2835	2852	2868	rBV	750017	1476831	4.93%	0.651%
29	9.732	2869	2887	2900	rBV	376584	751131	2.51%	0.331%
30	9.802	2901	2913	2916	rBV4	199330	311675	1.04%	0.137%
31	9.872	2930	2939	2959	rVB2	656992	1310899	4.38%	0.578%
32	9.971	2962	2976	2986	rBV2	483122	911048	3.04%	0.401%
33	10.076	3006	3015	3047	rVB4	355153	912990	3.05%	0.402%
34	10.207	3048	3064	3090	rBV5	364771	975155	3.26%	0.430%
35	10.320	3091	3106	3132	rBV2	789087	1651654	5.52%	0.728%
36	10.438	3133	3150	3162	rBV2	3455610	6961572	23.26%	3.067%

LSC Area Percent Report

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Sample : N1278-04
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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\82N011822W.M
Title : SW846 8260

37	10.620	3202	3218	3235	rVB9	909052	2294594	7.67%	1.011%
38	10.781	3266	3278	3293	rVB	942621	1744015	5.83%	0.768%
39	10.864	3293	3309	3329	rBV5	727832	1852803	6.19%	0.816%
40	11.130	3393	3408	3415	rBV3	170987	369283	1.23%	0.163%
41	11.258	3448	3456	3459	rVV3	300337	433848	1.45%	0.191%
42	11.291	3461	3468	3478	rVV	838495	1378611	4.61%	0.607%
43	11.344	3479	3488	3500	rVV	653692	1138249	3.80%	0.501%
44	11.449	3518	3527	3539	rVB5	207915	378227	1.26%	0.167%
45	11.548	3548	3564	3584	rVB10	242605	626864	2.09%	0.276%
46	11.744	3622	3637	3658	rBV	2208563	4088569	13.66%	1.801%
47	11.843	3659	3674	3694	rVV	3337351	6224348	20.80%	2.742%
48	11.948	3695	3713	3754	rVB	9629313	20451243	68.34%	9.010%
49	12.275	3813	3835	3845	rBV2	286848	887786	2.97%	0.391%
50	12.452	3891	3901	3911	rBV2	337065	550151	1.84%	0.242%
51	12.575	3936	3947	3961	rVB	932091	1522750	5.09%	0.671%
52	12.728	3991	4004	4024	rBV	1657785	2919349	9.75%	1.286%
53	12.916	4058	4074	4086	rBV	958739	1751989	5.85%	0.772%
54	12.980	4086	4098	4106	rBV	2544812	4333974	14.48%	1.909%
55	13.055	4118	4126	4141	rVB	1706255	2811859	9.40%	1.239%
56	13.219	4170	4187	4203	rBV	934639	1695384	5.66%	0.747%
57	13.318	4205	4224	4229	rBV7	150282	375642	1.26%	0.165%
58	13.367	4229	4242	4271	rVB	5021752	8432121	28.18%	3.715%
59	13.495	4275	4290	4301	rBV3	318626	631823	2.11%	0.278%
60	13.557	4302	4313	4324	rBV	530808	852034	2.85%	0.375%
61	13.672	4343	4356	4361	rBV	1869173	3093929	10.34%	1.363%
62	13.839	4406	4418	4420	rBV	323896	413820	1.38%	0.182%
63	13.871	4421	4430	4438	rBV3	1019212	1534365	5.13%	0.676%
64	13.997	4466	4477	4490	rVB4	278778	467623	1.56%	0.206%
65	14.067	4491	4503	4514	rBV	389070	623968	2.08%	0.275%
66	14.128	4515	4526	4531	rBV	496342	774076	2.59%	0.341%
67	14.214	4548	4558	4570	rVB	717051	1046594	3.50%	0.461%
68	14.278	4571	4582	4588	rBV	239193	383059	1.28%	0.169%
69	14.324	4588	4599	4612	rVB2	1271215	2115586	7.07%	0.932%
70	14.458	4638	4649	4660	rBV2	304770	495785	1.66%	0.218%
71	14.539	4661	4679	4685	rVV2	324654	575362	1.92%	0.253%
72	14.576	4686	4693	4703	rVB	513430	752676	2.51%	0.332%
73	14.807	4770	4779	4790	rBV3	301703	529982	1.77%	0.233%
74	14.919	4805	4821	4838	rBV3	1153036	2740140	9.16%	1.207%
75	15.311	4952	4967	4983	rVB2	406308	964398	3.22%	0.425%
76	15.501	5026	5038	5054	rVB	589217	1138136	3.80%	0.501%

LSC Area Percent Report

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Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 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\82N011822W.M
Title : SW846 8260

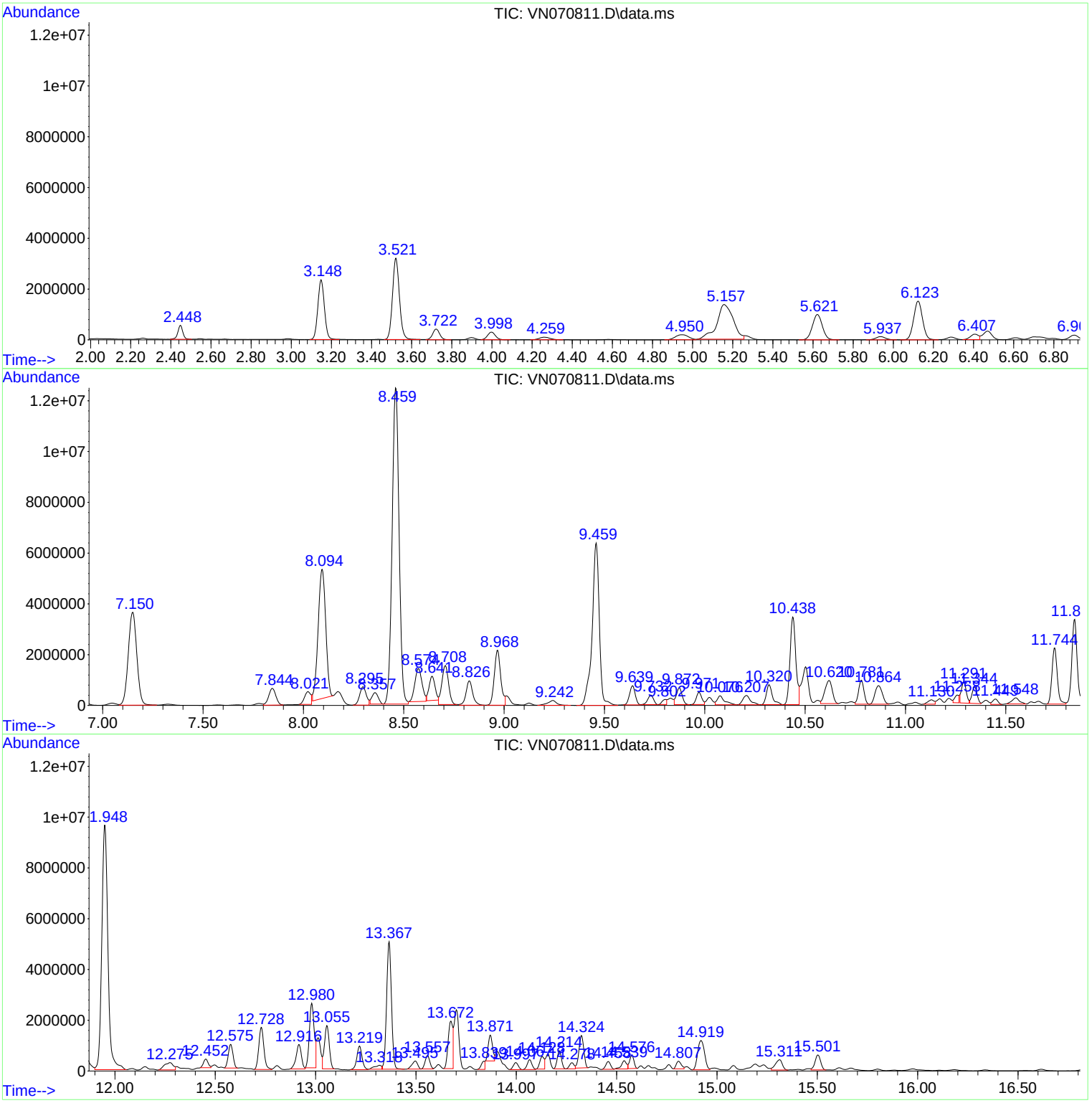
Sum of corrected areas: 226977572

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
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ClientSampleId :
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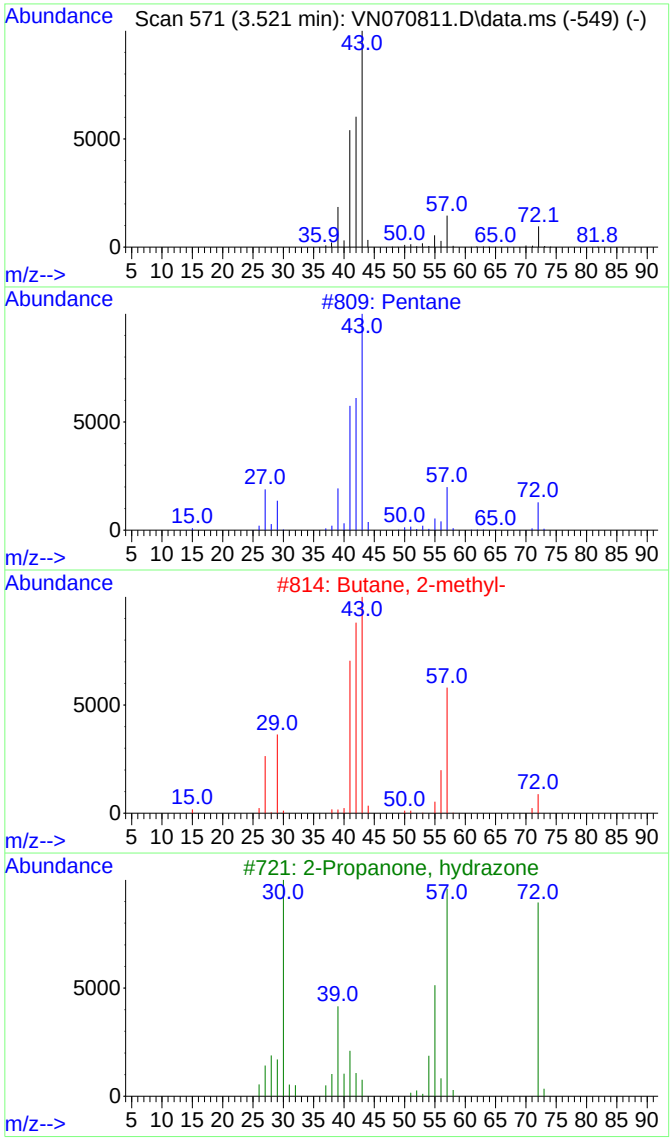
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

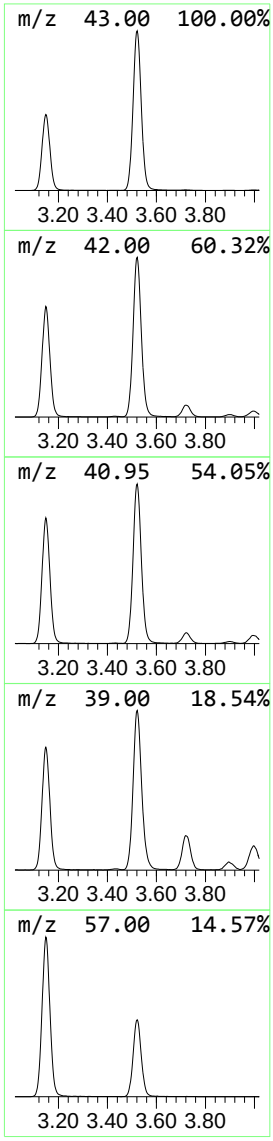
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Pentane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.521	28.91 ug/l	7711730	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	50
3	2-Propanone, hydrazone	72	C3H8N2	005281-20-9	9
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	9
5	Isobutyl nitrite	103	C4H9NO2	000542-56-3	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
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Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

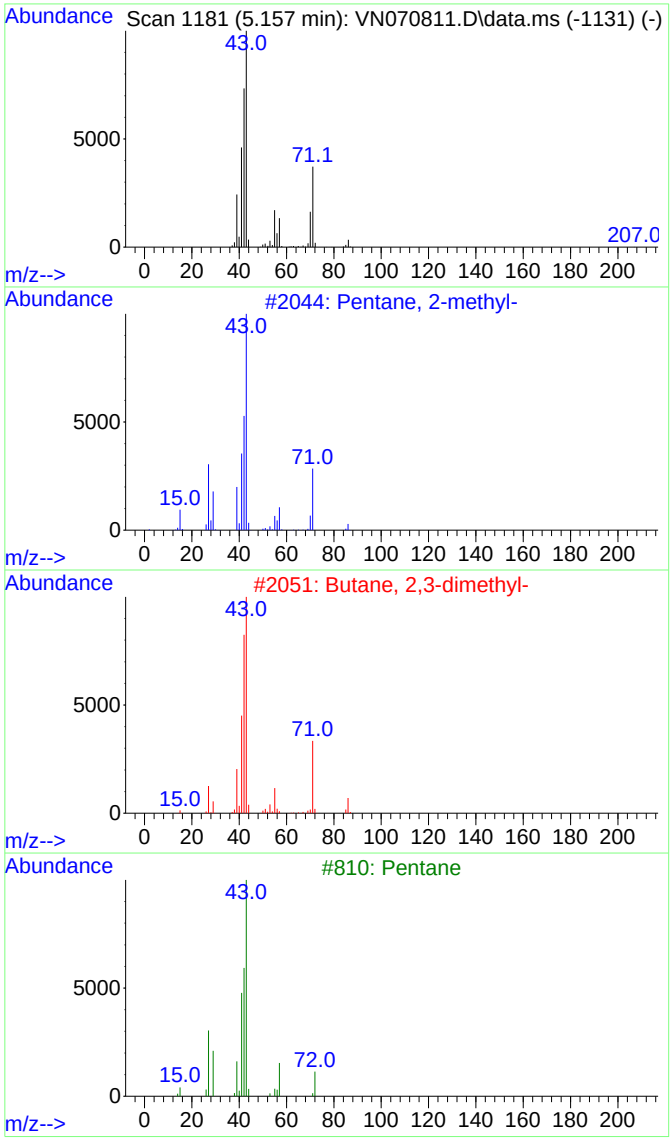
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

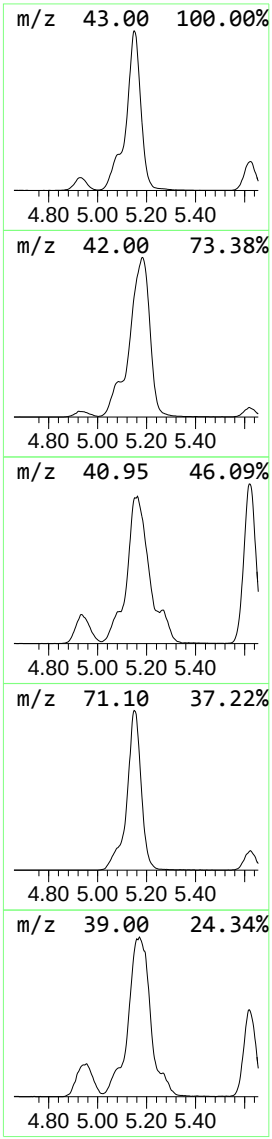
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.157	29.10 ug/l	7764340	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
3	Pentane	72	C5H12	000109-66-0	38
4	Cyclobutane, methyl-	70	C5H10	000598-61-8	25
5	1-Pentene	70	C5H10	000109-67-1	25





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

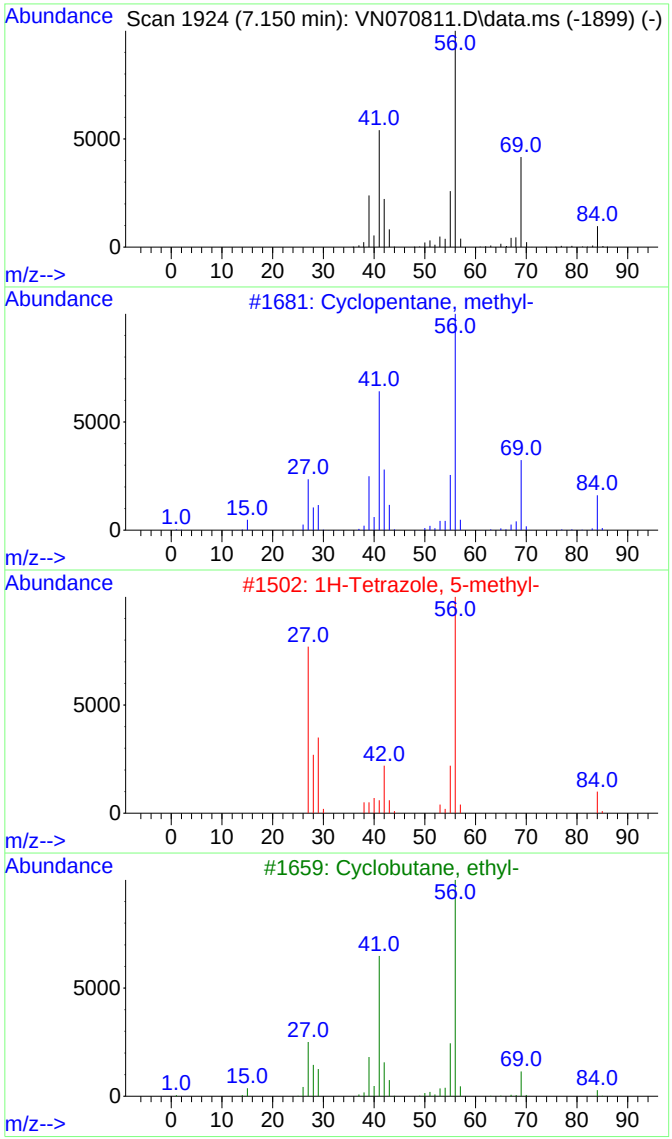
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

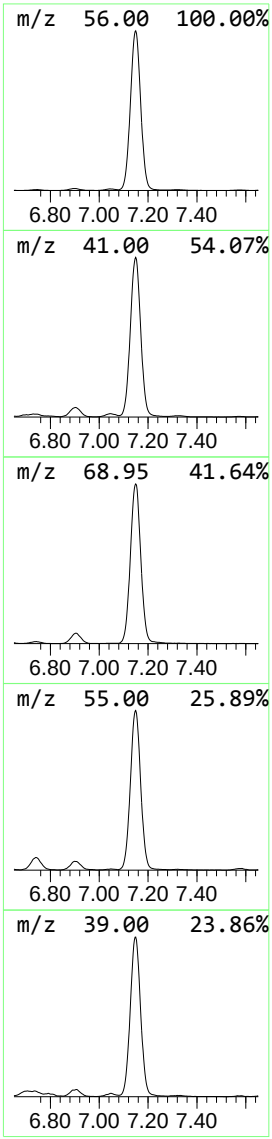
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclopentane, methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.150	40.48 ug/l	10799600	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4	Cyclohexane	84	C6H12	000110-82-7	56
5	Cyclobutane, butyl-	112	C8H16	013152-44-8	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

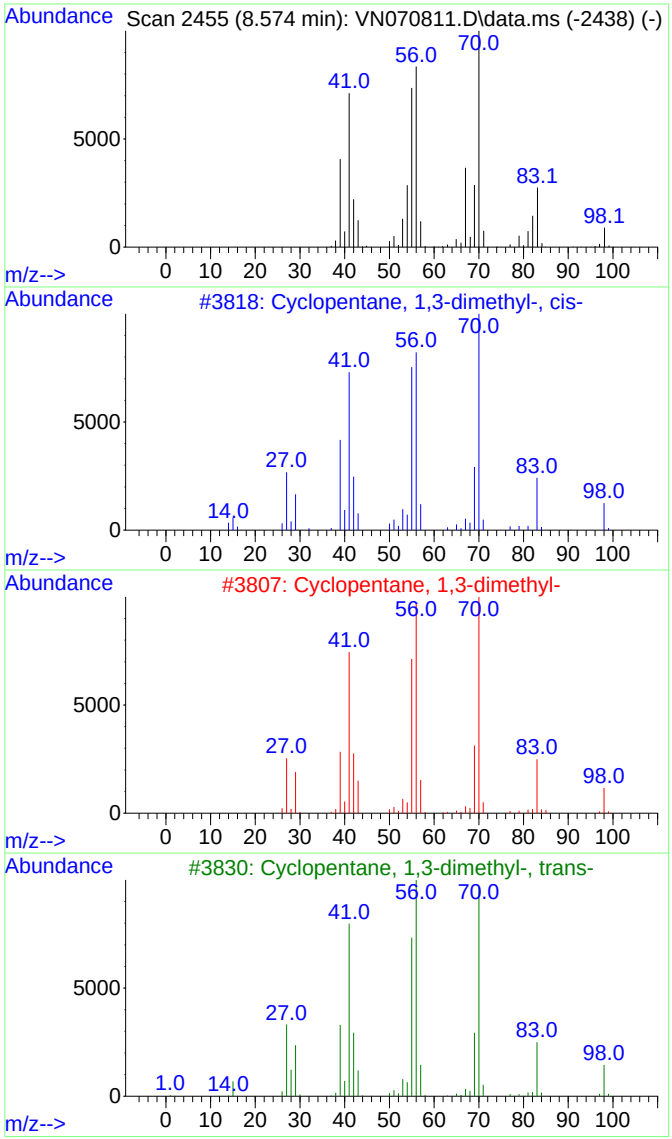
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

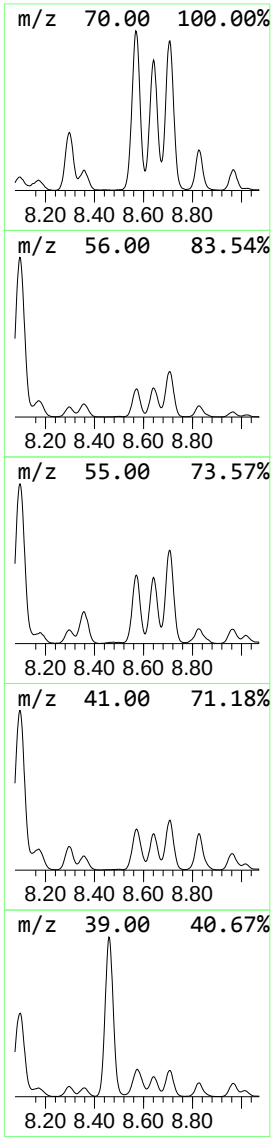
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclopentane, 1,3-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.574	33.74 ug/l	3254960	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	93
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	92
3	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	70
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	64
5	1-Hexene, 3-methyl-	98	C7H14	003404-61-3	38





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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
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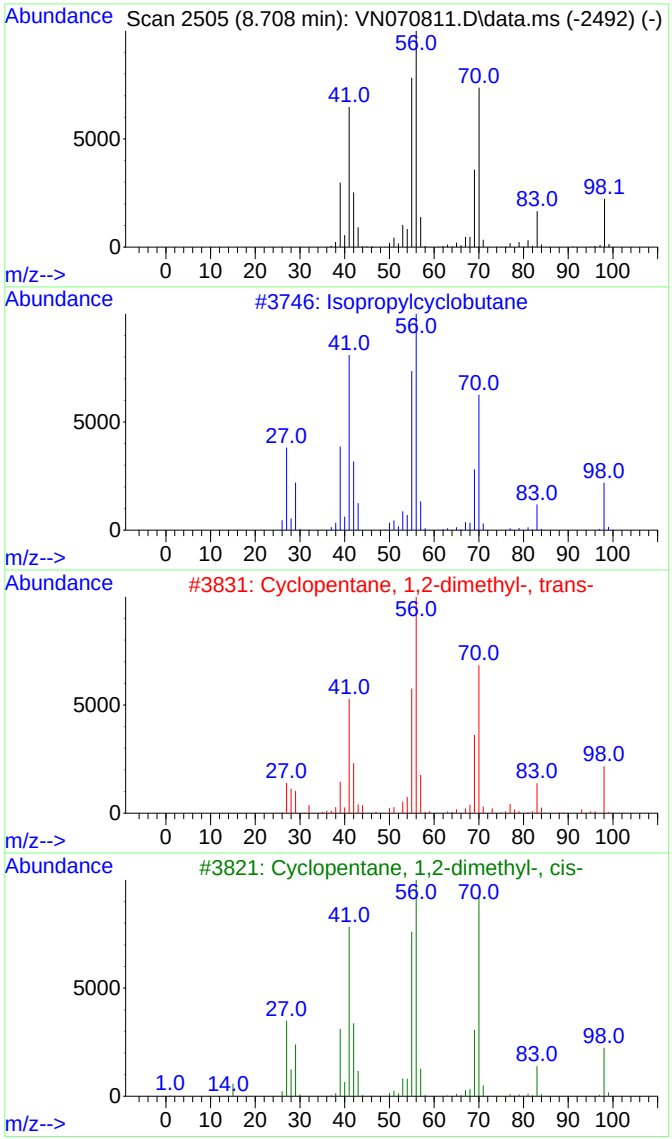
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

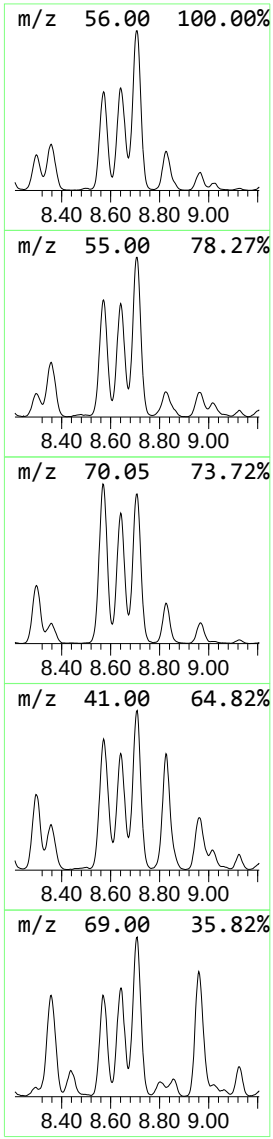
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Isopropylcyclobutane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.708	36.78 ug/l	3548660	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropylcyclobutane	98	C7H14	000872-56-0	94
2	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
3	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	91
4	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
5	1-Heptene	98	C7H14	000592-76-7	83





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Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

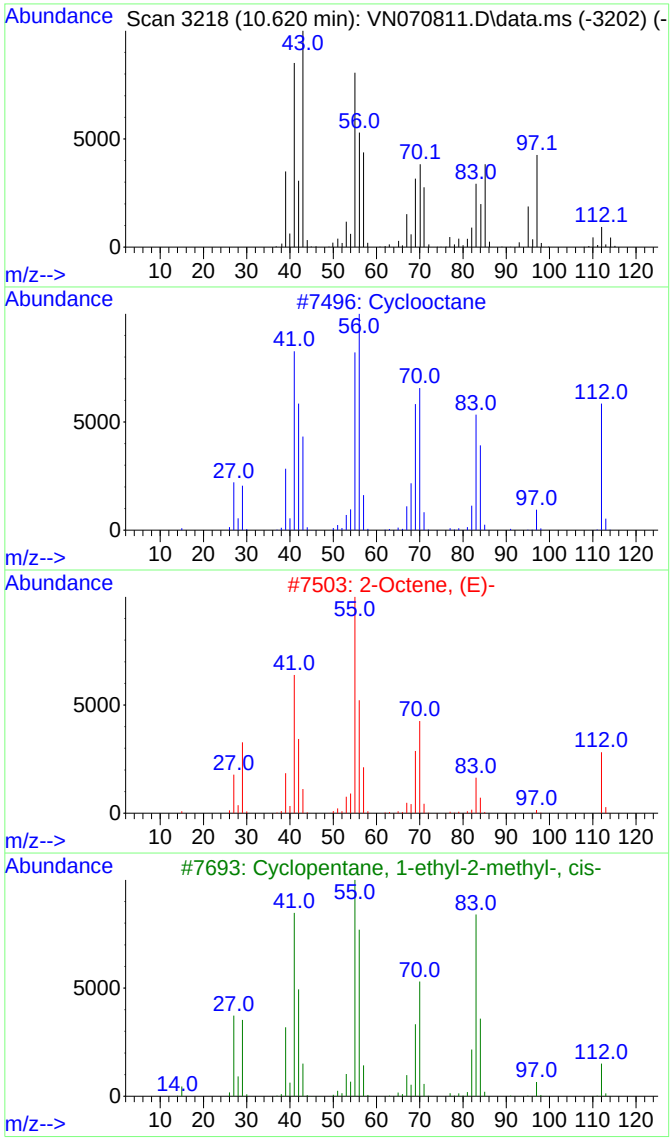
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

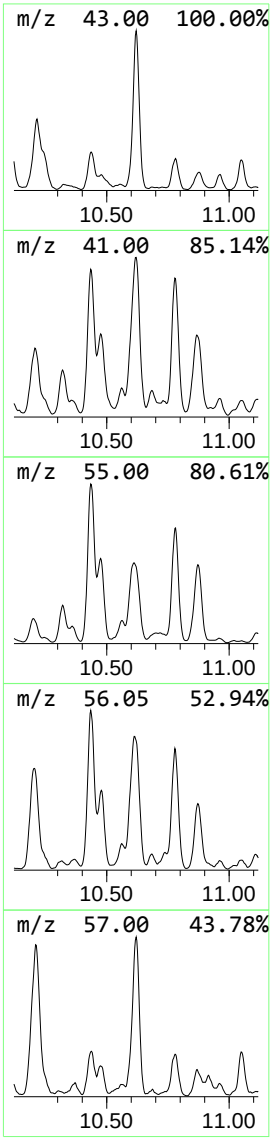
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 unknown10.620 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.620	28.06 ug/l	2294590	Chlorobenzene-d5	11.744

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclooctane	112	C8H16	000292-64-8	49
2	2-Octene, (E)-	112	C8H16	013389-42-9	38
3	Cyclopentane, 1-ethyl-2-methyl-,...	112	C8H16	000930-89-2	38
4	Cyclopentane, 1-ethyl-2-methyl-	112	C8H16	003726-46-3	38
5	2-Octene	112	C8H16	000111-67-1	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

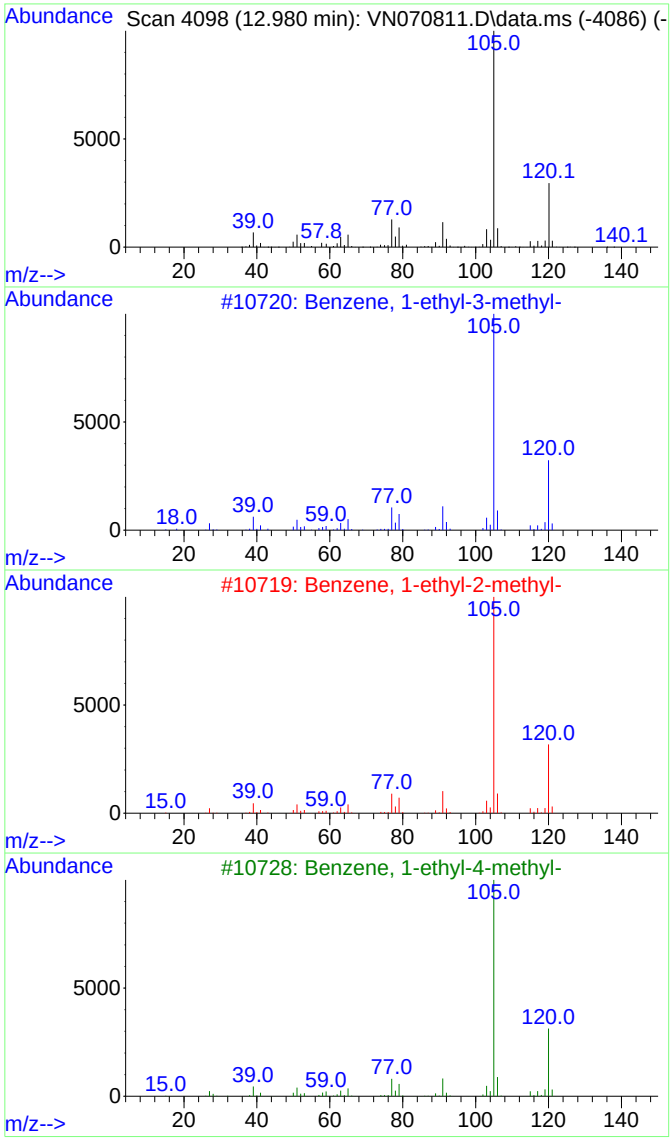
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

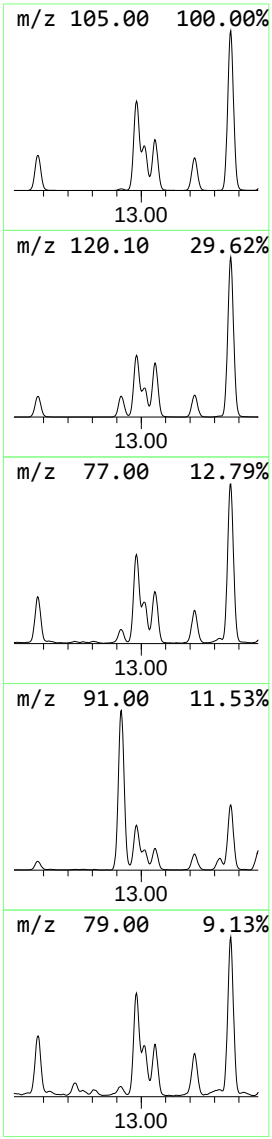
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	70.04 ug/l	4333970	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
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Operator : JC/MD
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ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

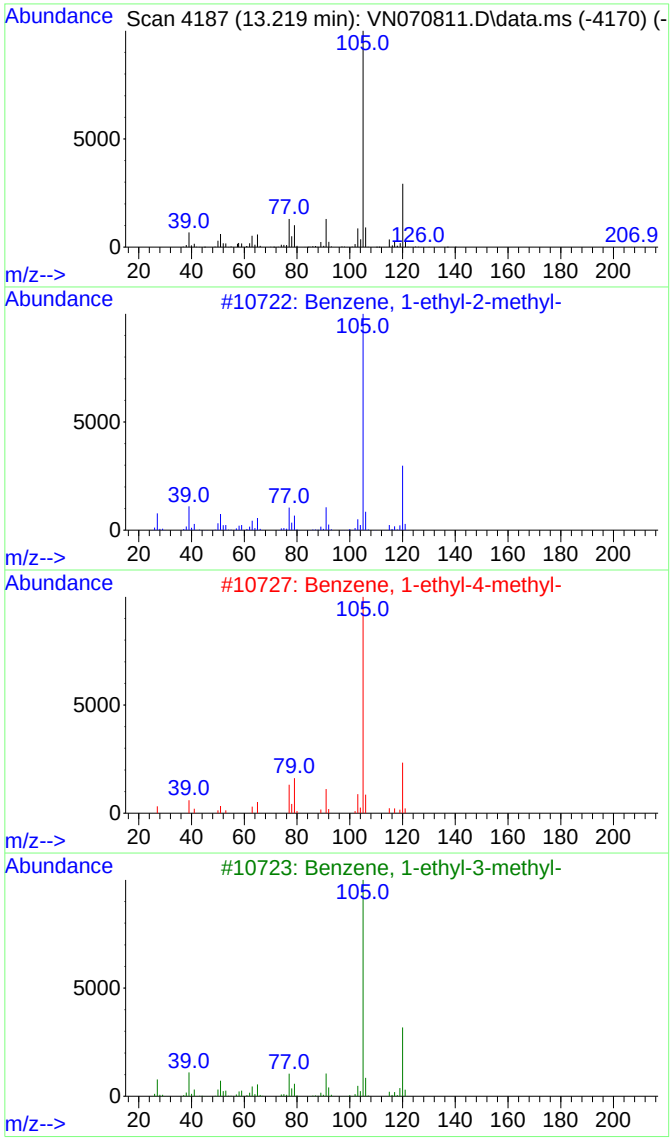
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

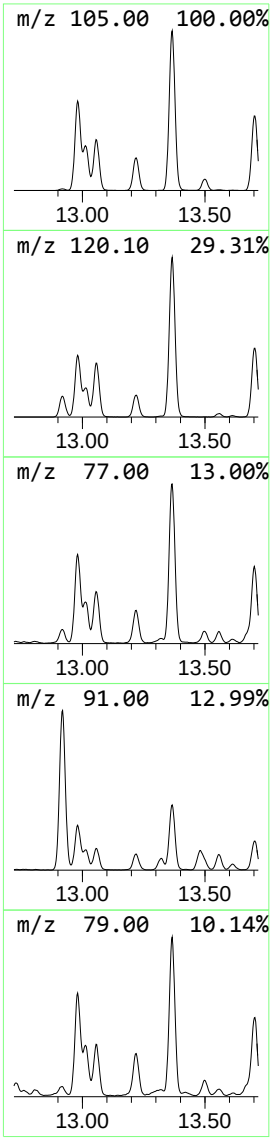
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.219	27.40 ug/l	1695380	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	95
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

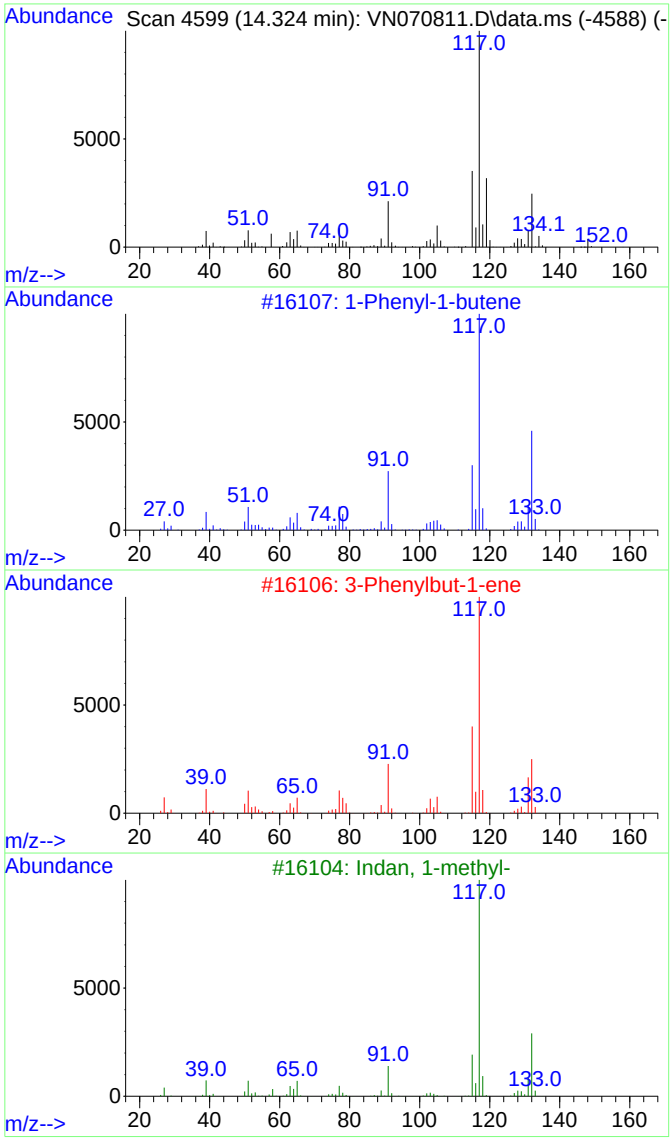
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

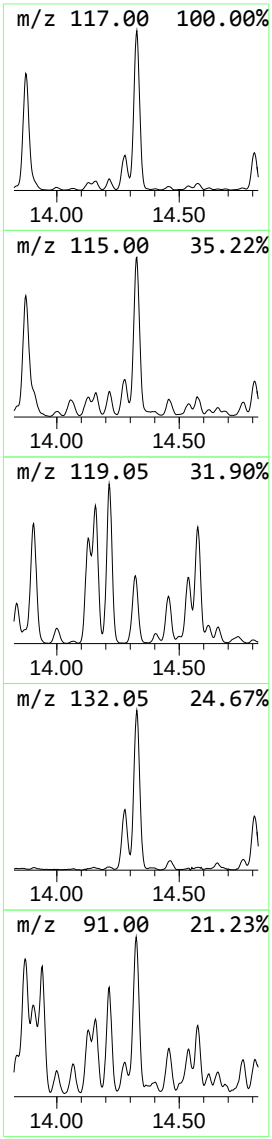
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 1-Phenyl-1-butene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.324	34.19 ug/l	2115590	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	81
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Indan, 1-methyl-	132	C10H12	000767-58-8	76
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	72
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

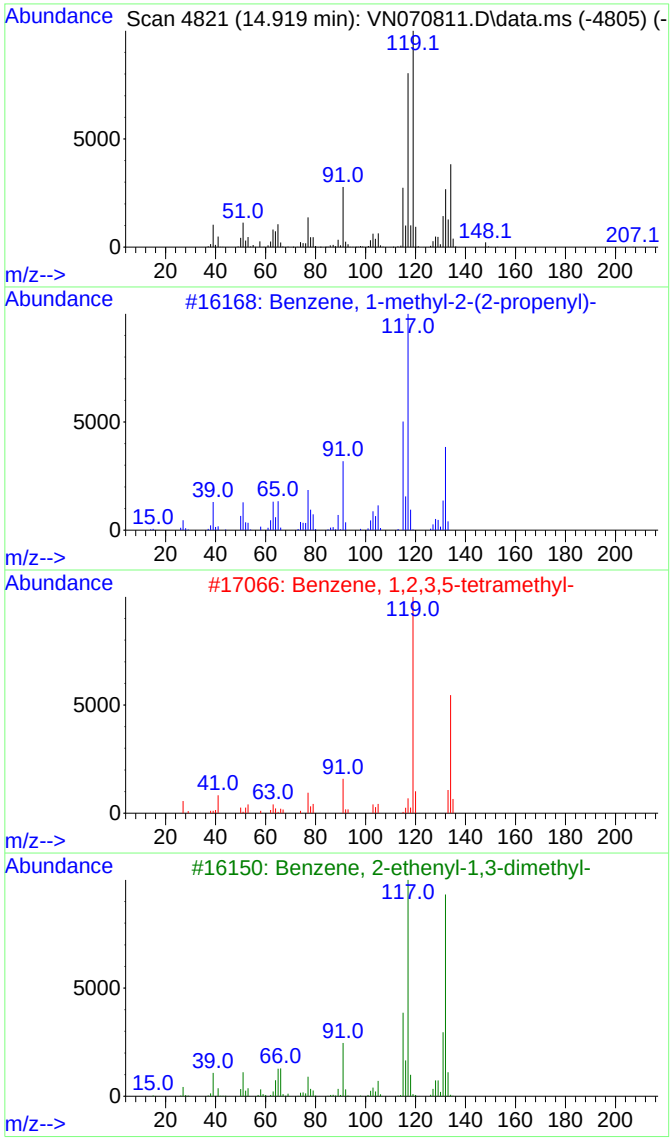
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.M
Quant Title : SW846 8260

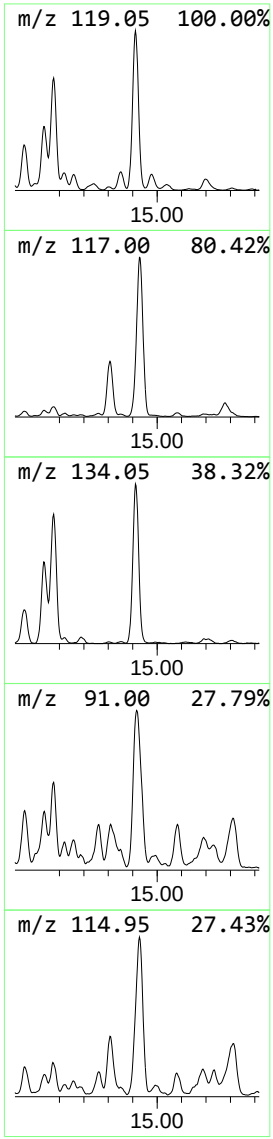
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-methyl-2-(2-prop... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.919	44.28 ug/l	2740140	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50
3	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	50
5	p-Cymene	134	C10H14	000099-87-6	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN012722\
Data File : VN070811.D
Acq On : 27 Jan 2022 17:21
Operator : JC/MD
Sample : N1278-04
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 14 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N011822W.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
Pentane	3.521	28.9	ug/l	7711730	1	8.086	13339700	50.0
Pentane, 2-methyl-	5.157	29.1	ug/l	7764340	1	8.086	13339700	50.0
Cyclopentane, m...	7.150	40.5	ug/l	10799600	1	8.086	13339700	50.0
Cyclopentane, 1...	8.574	33.7	ug/l	3254960	2	8.968	4823540	50.0
Isopropylcyclob...	8.708	36.8	ug/l	3548660	2	8.968	4823540	50.0
unknown10.620	10.620	28.1	ug/l	2294590	3	11.744	4088570	50.0
Benzene, 1-ethy...	12.980	70.0	ug/l	4333970	4	13.672	3093930	50.0
Benzene, 1-ethy...	13.219	27.4	ug/l	1695380	4	13.672	3093930	50.0
1-Phenyl-1-butene	14.324	34.2	ug/l	2115590	4	13.672	3093930	50.0
Benzene, 1-meth...	14.919	44.3	ug/l	2740140	4	13.672	3093930	50.0