LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

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\82N120221W.M
Title : SW846 8260

Signal : TIC: VN070102.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.613	1326	1351	1376	rBV4	70399	241830	1.85%	0.098%
2	6.117	1513	1539	1562	rBV3	94844	293344	2.24%	0.119%
3	7.042	1861	1884	1905	rBV4	142455	425375	3.25%	0.172%
4	7.144	1906	1922	1942	rVB3	63594	183293	1.40%	0.074%
5	8.021	2227	2249	2257	rBV	917698	2195340	16.79%	0.888%
6	8.080	2259	2271	2290	rVV2	2155603	5662028	43.30%	2.291%
7	8.174	2292	2306	2330	rVB2	517137	1314235	10.05%	0.532%
8	8.295	2331	2351	2384	rBV2	1029232	2406323	18.40%	0.974%
9	8.437	2385	2404	2433	rBV	1120220	2593849	19.83%	1.049%
10	8.614	2434	2470	2494	rBV	2869379	8007997	61.23%	3.240%
11	8.708	2496	2505	2526	rVB3	159399	408463	3.12%	0.165%
12	8.826	2527	2549	2577	rVB2	823696	1838008	14.05%	0.744%
13	8.965	2580	2601	2628	rBV	3323144	7422098	56.75%	3.003%
14	9.123	2649	2660	2674	rVB3	192668	356960	2.73%	0.144%
15	9.199	2674	2688	2698	rBV3	193316	390967	2.99%	0.158%
16	9.456	2752	2784	2795	rBV2	1971557	4780218	36.55%	1.934%
17	9.510	2796	2804	2821	rVV	1132568	2197342	16.80%	0.889%
18	9.593	2821	2835	2843	rVV	307459	618672	4.73%	0.250%
19	9.638	2844	2852	2865	rVV4	316630	617054	4.72%	0.250%
20	9.732	2866	2887	2903	rVV5	362593	971984	7.43%	0.393%
21	9.883	2927	2943	2965	rBV	3285093	6968529	53.29%	2.819%
22	10.014	2966	2992	3005	rBV3	4613078	10625219	81.25%	4.299%
23	10.078	3006	3016	3025	rBV	2033075	3556318	27.19%	1.439%
24	10.215	3047	3067	3093	rBV	2560684	6060916	46.35%	2.452%
25	10.317	3094	3105	3111	rBV6	215732	395934	3.03%	0.160%
26	10.371	3112	3125	3137	rVV2	2482099	4547304	34.77%	1.840%
27	10.440	3137	3151	3174	rVV	5614175	10943765	83.68%	4.428%
28	10.556	3175	3194	3203	rBV10	435007	1333183	10.19%	0.539%
29	10.623	3205	3219	3233	rVV3	2153369	4570817	34.95%	1.849%
30	10.682	3233	3241	3251	rVB3	508662	774115	5.92%	0.313%
31	10.789	3268	3281	3283	rBV4	456198	701952	5.37%	0.284%
32	10.864	3296	3309	3332	rVB7	1262258	3869965	29.59%	1.566%
33	10.963	3336	3346	3358	rVB2	712554	1305185	9.98%	0.528%
34	11.052	3359	3379	3391	rBV3	784775	1714509	13.11%	0.694%
35	11.151	3392	3416	3433	rBV3	1223178	3628612	27.75%	1.468%
36	11.218	3434	3441	3446	rBV6	300232	422241	3.23%	0.171%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

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\82N120221W.M
Title : SW846 8260

37	11.256	3449	3455	3463	rVB4	404530	528167	4.04%	0.214%
38	11.344	3482	3488	3496	rBV2	315238	452741	3.46%	0.183%
39	11.454	3521	3529	3538	rBV2	693807	1013919	7.75%	0.410%
40	11.527	3545	3556	3581	rVB4	3515137	8036463	61.45%	3.251%
41	11.626	3582	3593	3615	rBV	2695427	4907068	37.52%	1.985%
42	11.744	3623	3637	3653	rBV	6024956	10709821	81.89%	4.333%
43	12.071	3753	3759	3772	rVB6	332756	504403	3.86%	0.204%
44	12.251	3815	3826	3838	rBV4	1002682	1938256	14.82%	0.784%
45	12.578	3937	3948	3959	rBV	3978760	6722710	51.41%	2.720%
46	12.728	3995	4004	4014	rVB2	3521379	5331417	40.77%	2.157%
47	12.919	4063	4075	4096	rVB	6579193	11082203	84.74%	4.484%
48	13.055	4117	4126	4141	rVB4	1570635	2982483	22.81%	1.207%
49	13.283	4206	4211	4218	rBV3	397399	466922	3.57%	0.189%
50	13.412	4253	4259	4271	rVB4	497901	787367	6.02%	0.319%
51	13.605	4326	4331	4337	rBV2	414073	402053	3.07%	0.163%
52	13.672	4347	4356	4364	rBV	4863381	8050251	61.56%	3.257%
53	13.836	4406	4417	4424	rBV	3083733	4939852	37.77%	1.999%
54	13.876	4425	4432	4438	rVV	1780235	2446406	18.71%	0.990%
55	13.999	4468	4478	4487	rBV4	954204	1545323	11.82%	0.625%
56	14.069	4496	4504	4516	rVB	1296191	1957683	14.97%	0.792%
57	14.131	4518	4527	4543	rVB	1147971	2023726	15.47%	0.819%
58	14.214	4546	4558	4571	rBV	8214651	13077753	100.00%	5.291%
59	14.324	4589	4599	4612	rVB4	3578203	6082638	46.51%	2.461%
60	14.458	4640	4649	4659	rBV2	1045088	1709318	13.07%	0.692%
61	14.538	4670	4679	4690	rVB	3937620	6080653	46.50%	2.460%
62	14.619	4701	4709	4718	rBV2	1652832	2356834	18.02%	0.954%
63	14.662	4719	4725	4741	rVB7	717152	1124890	8.60%	0.455%
64	14.809	4769	4780	4791	rBV2	2912097	5057179	38.67%	2.046%
65	14.927	4807	4824	4836	rBV3	5157558	12107834	92.58%	4.899%
66	14.978	4838	4843	4860	rVB	1139563	1904463	14.56%	0.771%
67	15.083	4875	4882	4898	rVB2	490952	796955	6.09%	0.322%
68	15.193	4913	4923	4936	rBV5	1674402	3033521	23.20%	1.227%
69	15.308	4954	4966	4982	rVB3	1628650	3679023	28.13%	1.489%
70	15.464	5016	5024	5028	rBV5	278434	400708	3.06%	0.162%
71	15.501	5031	5038	5052	rVB	1069554	1815972	13.89%	0.735%
72	15.611	5070	5079	5089	rBV3	546615	911432	6.97%	0.369%
73	15.802	5137	5150	5168	rBV	944800	1928454	14.75%	0.780%
74	16.614	5439	5453	5497	rVB	1609242	3921346	29.98%	1.587%

Sum of corrected areas: 247162151

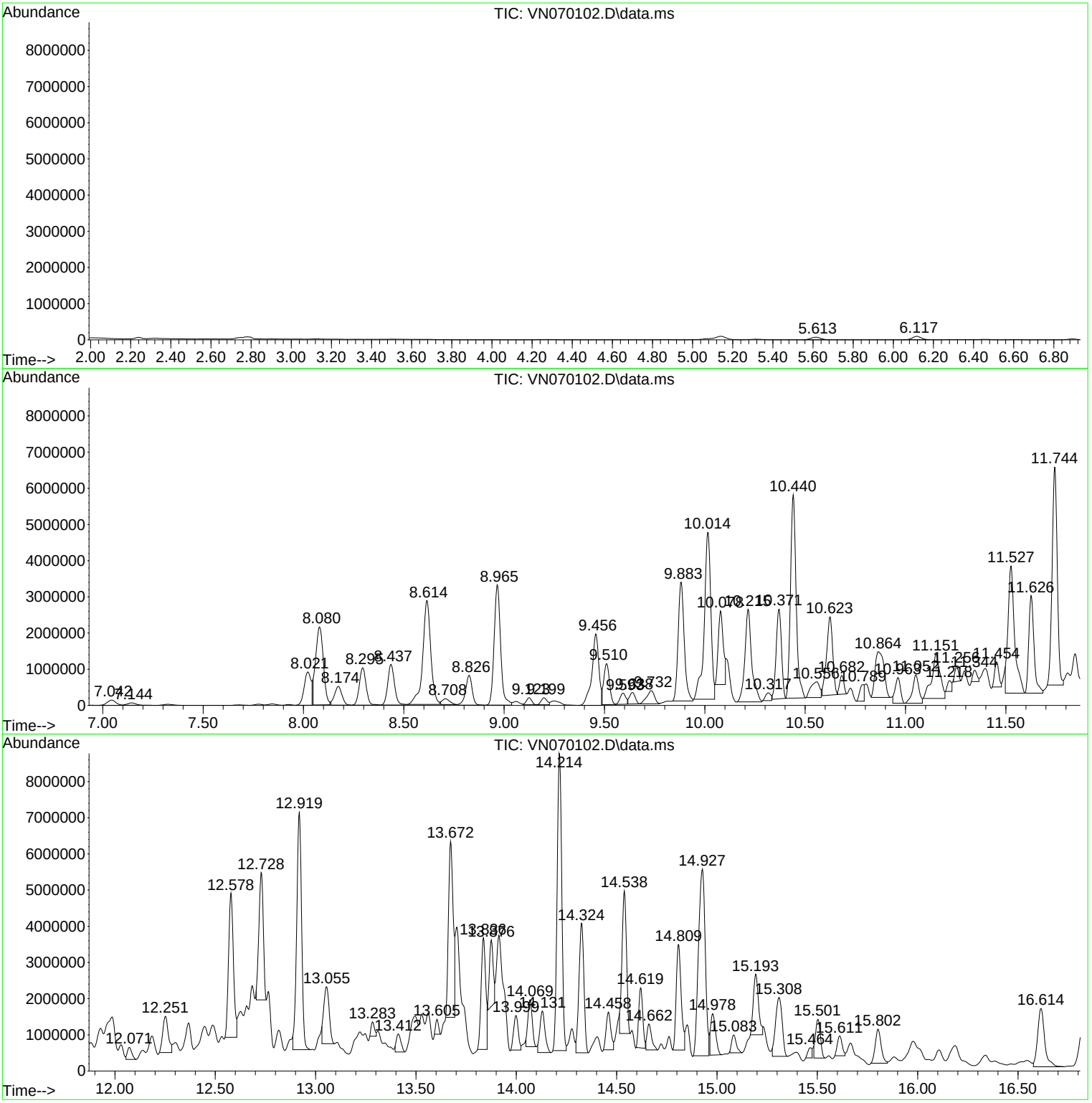
Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

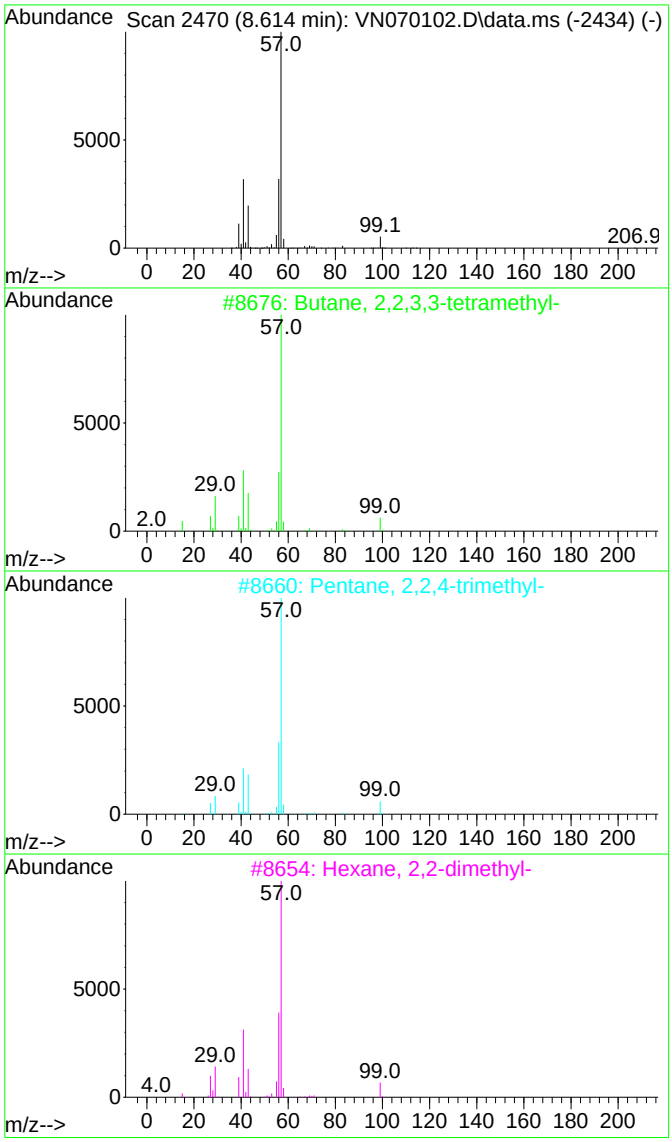
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

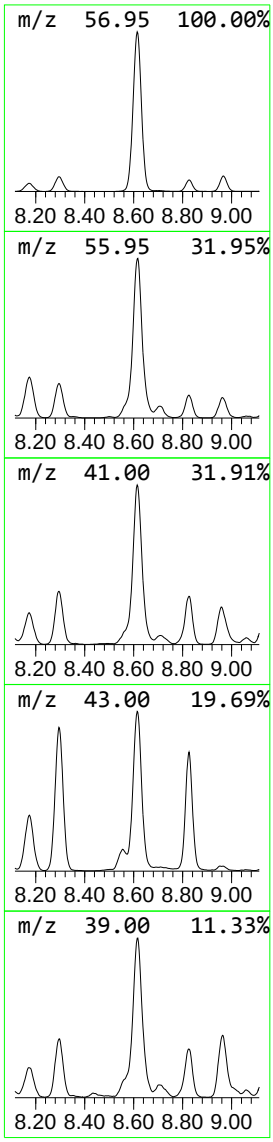
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Butane, 2,2,3,3-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.614	53.95 ug/l	8008000	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,2,3,3-tetramethyl-	114	C8H18	000594-82-1	83
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	78
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	78
4	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	64
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

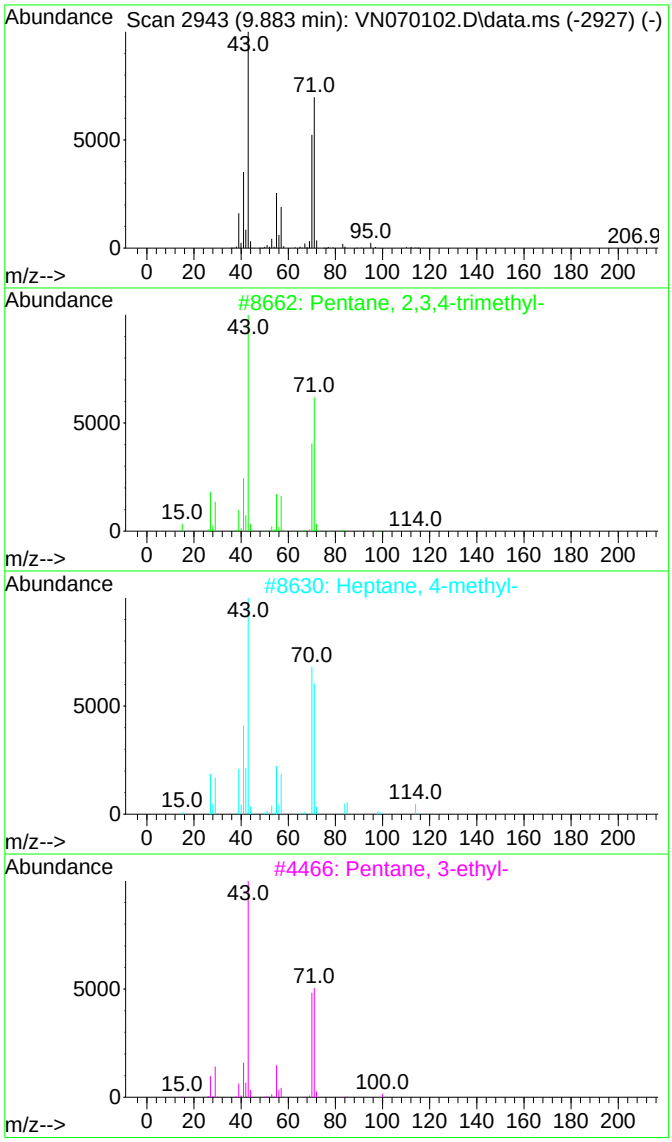
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

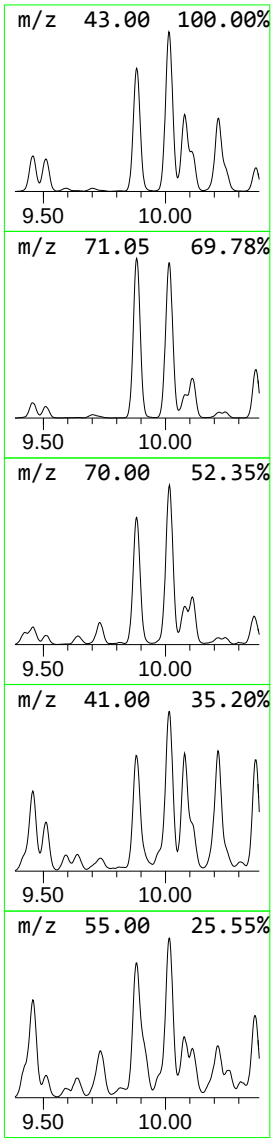
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Pentane, 2,3,4-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.883	46.94 ug/l	6968530	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	91
2	Heptane, 4-methyl-	114	C8H18	000589-53-7	83
3	Pentane, 3-ethyl-	100	C7H16	000617-78-7	56
4	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	56
5	Hexane, 3-methyl-	100	C7H16	000589-34-4	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

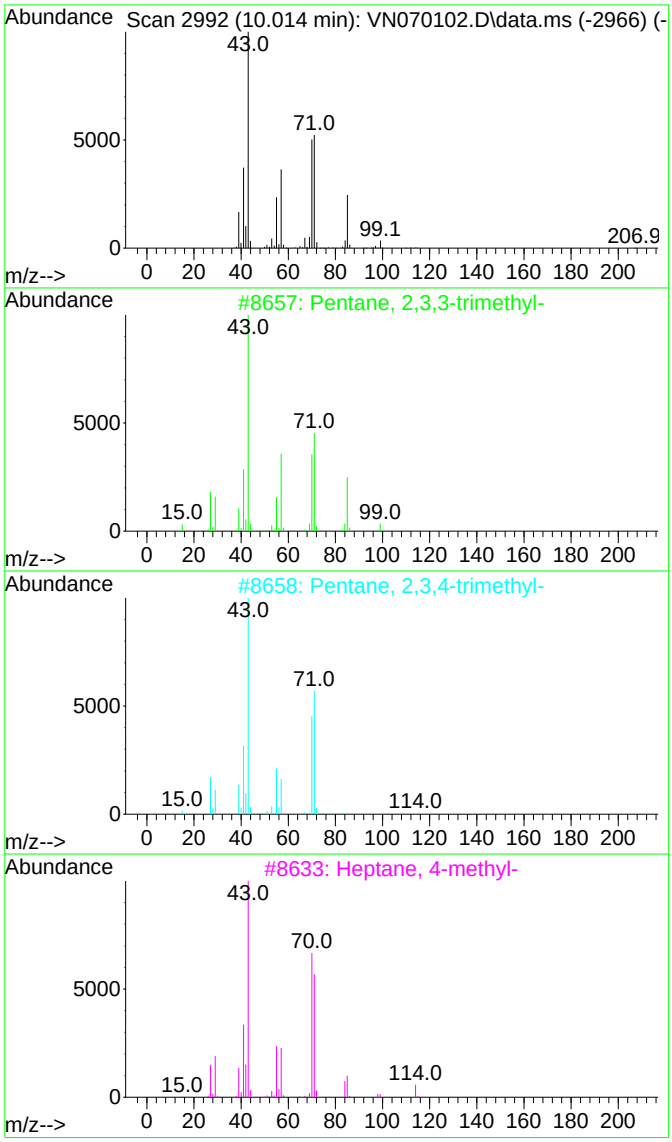
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

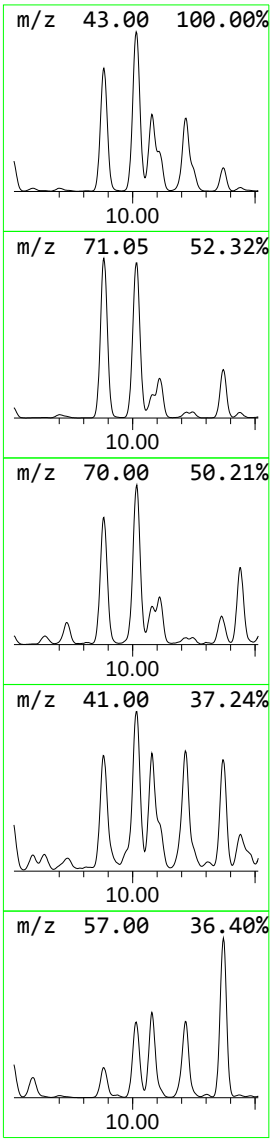
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.014	71.58 ug/l	10625200	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	68
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	53
5	Pentane, 3-ethyl-2-methyl-	114	C8H18	000609-26-7	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

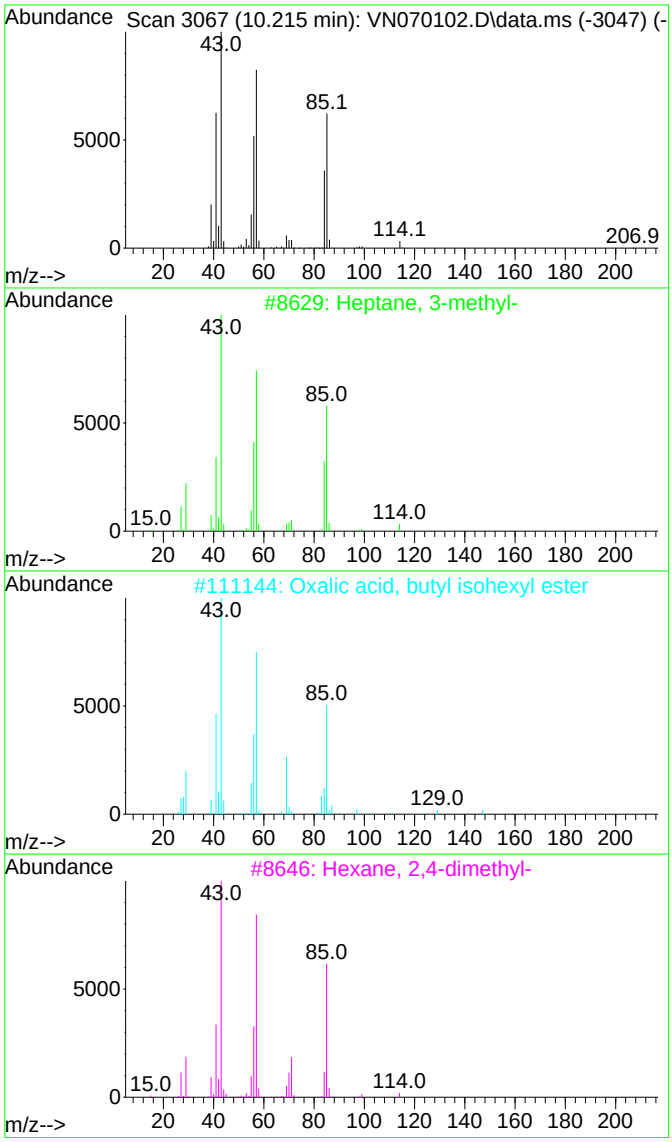
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

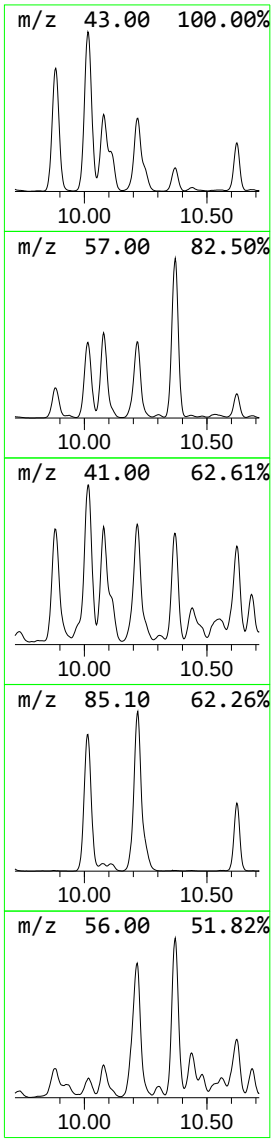
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Heptane, 3-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.215	40.83 ug/l	6060920	1,4-Difluorobenzene	8.968

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2	Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	78
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	64
4	Octane, 3,4-dimethyl-	142	C10H22	015869-92-8	59
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

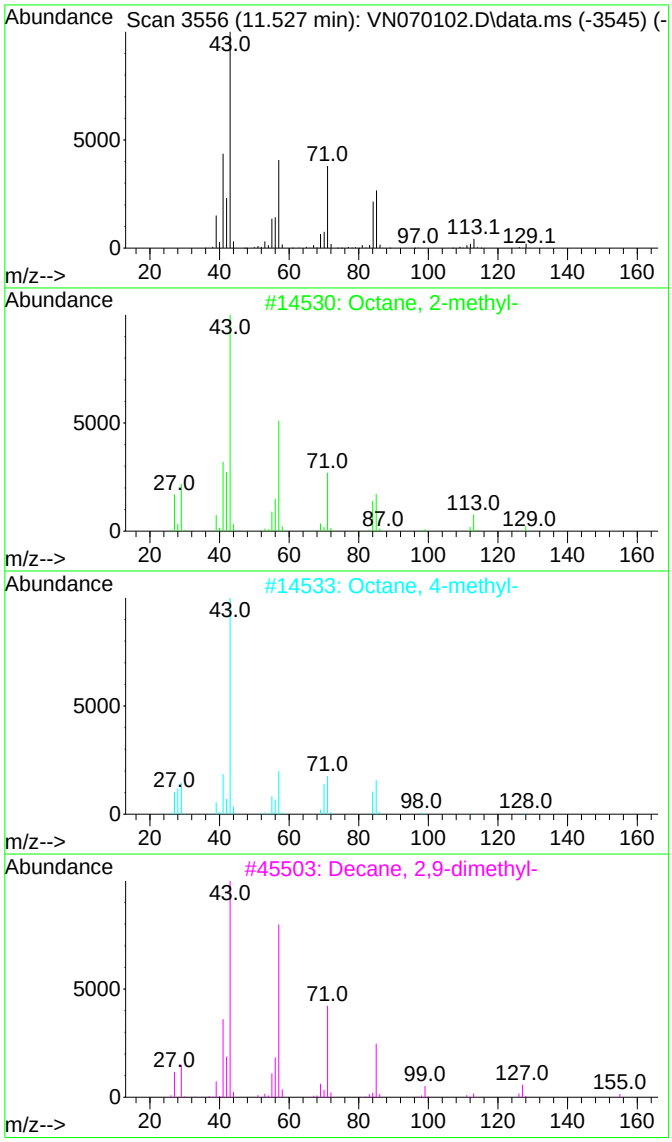
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

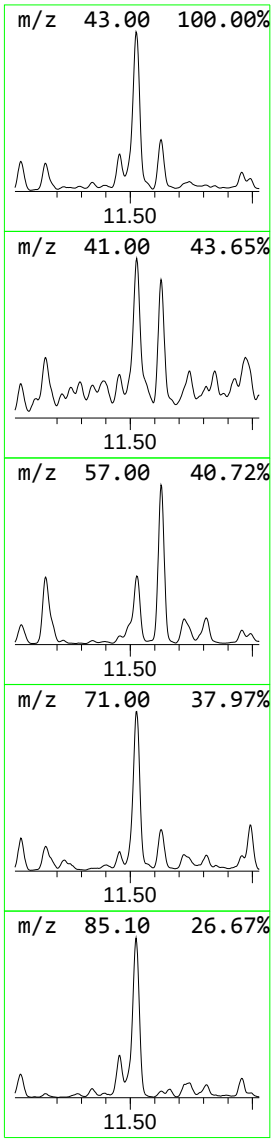
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Octane, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.527	37.52 ug/l	8036460	Chlorobenzene-d5	11.744

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2-methyl-	128	C9H20	003221-61-2	83
2	Octane, 4-methyl-	128	C9H20	002216-34-4	59
3	Decane, 2,9-dimethyl-	170	C12H26	001002-17-1	59
4	Hexane, 3-ethyl-	114	C8H18	000619-99-8	53
5	Decane, 2,4-dimethyl-	170	C12H26	002801-84-5	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

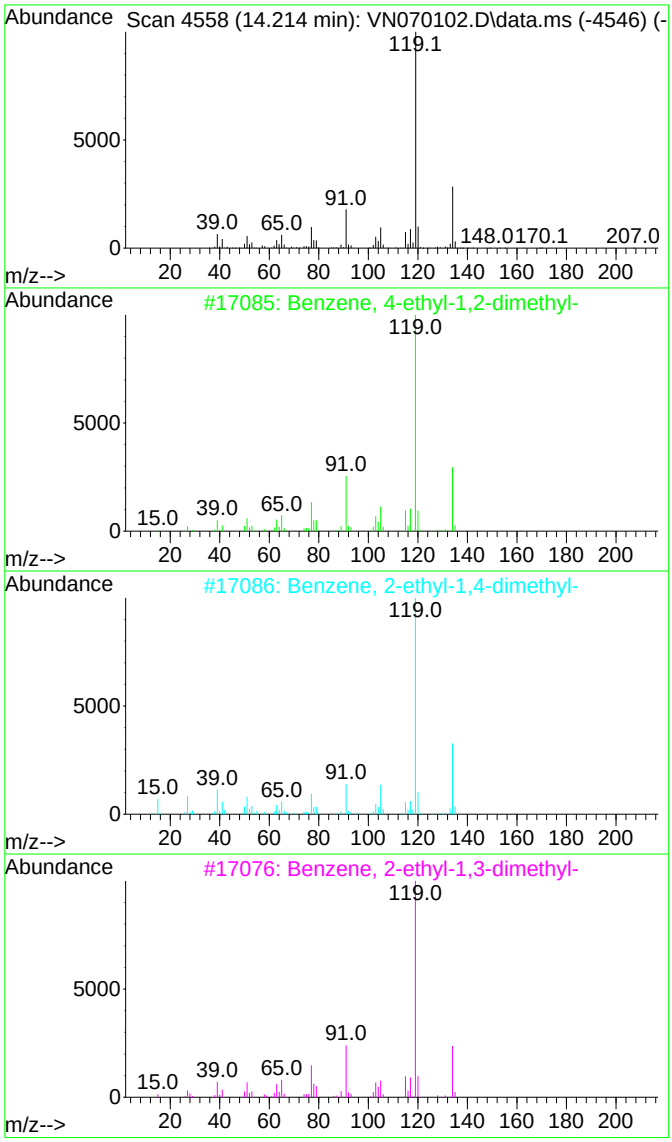
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

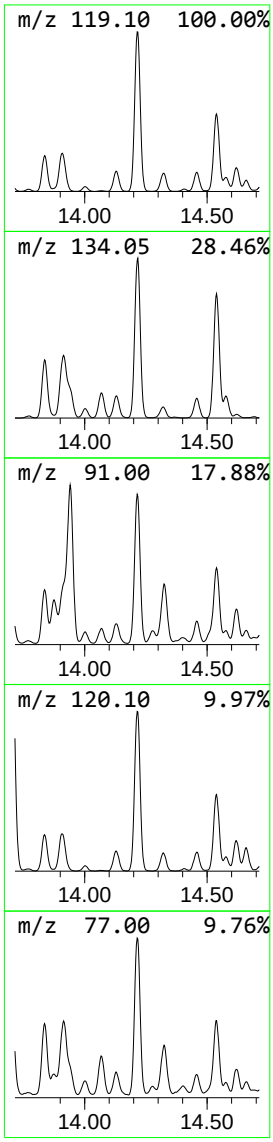
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.214	81.23 ug/l	13077800	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

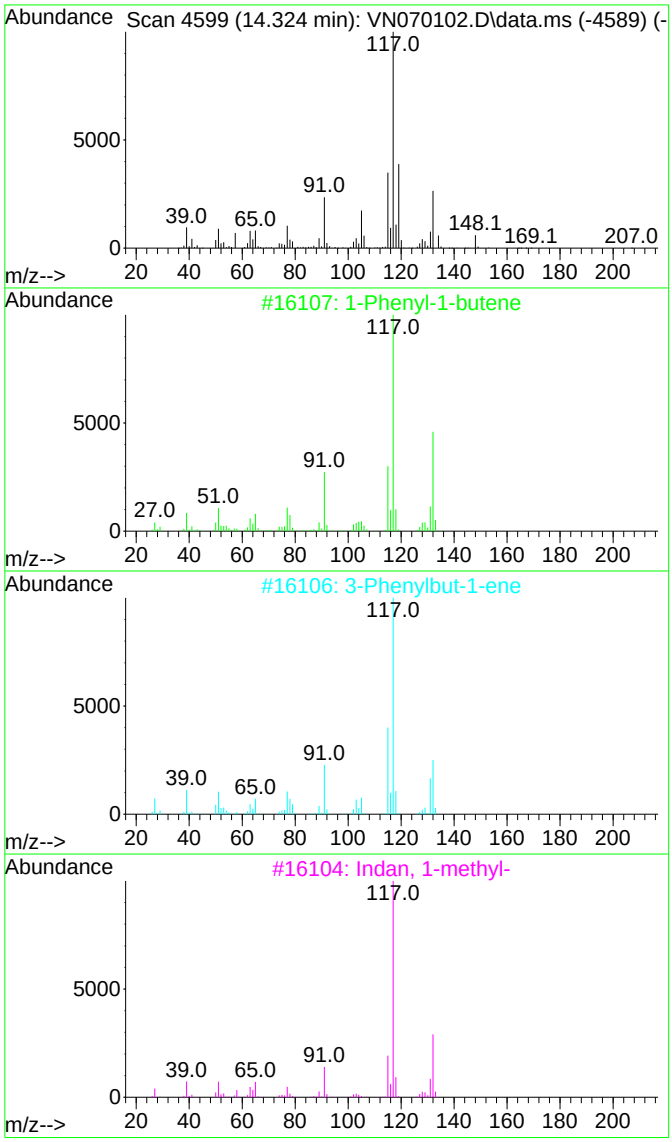
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

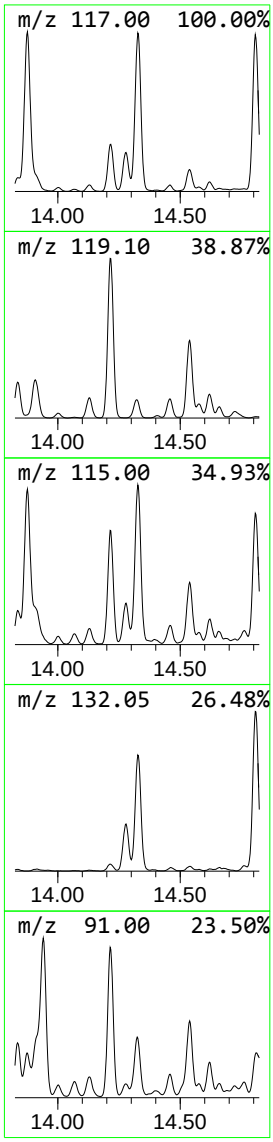
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1-Phenyl-1-butene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.324	37.78 ug/l	6082640	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	86
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
3	Indan, 1-methyl-	132	C10H12	000767-58-8	70
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	70
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	62





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

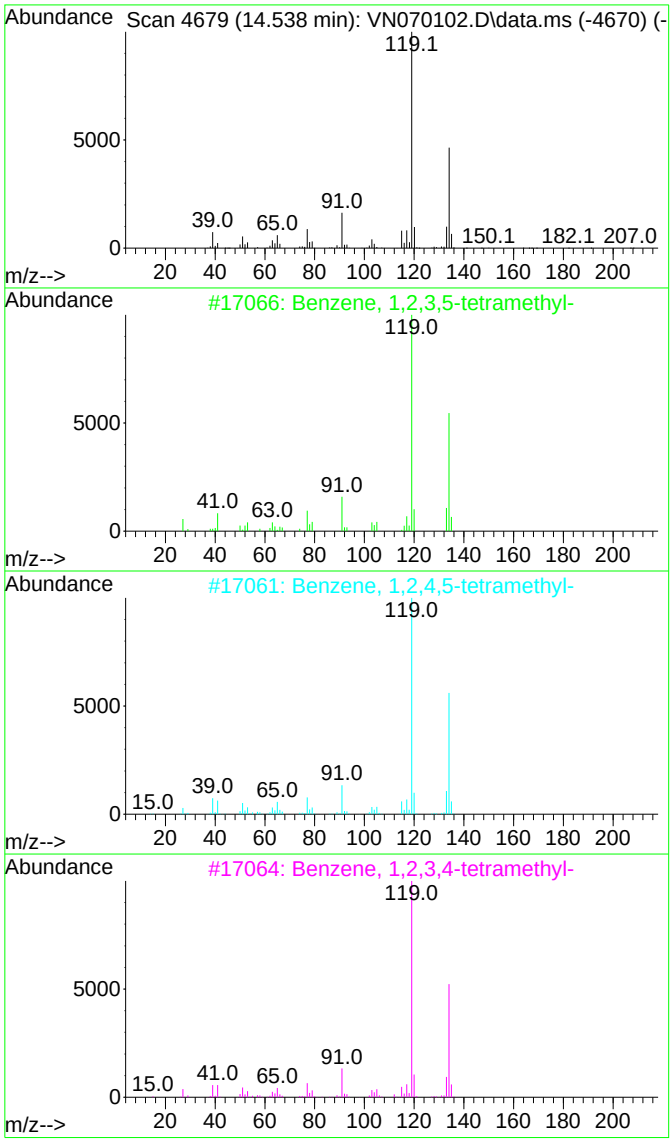
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

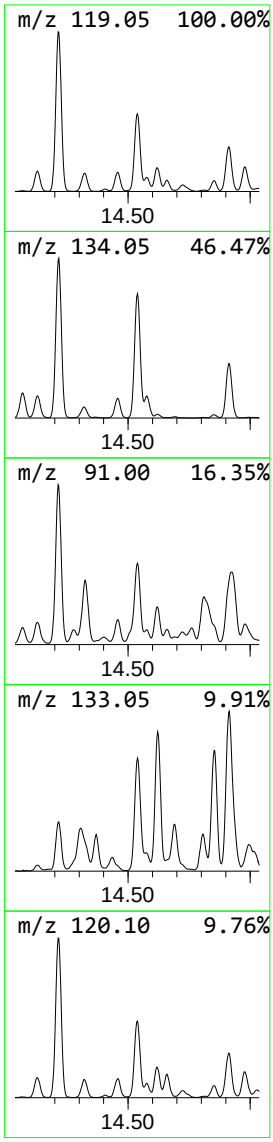
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.538	37.77 ug/l	6080650	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

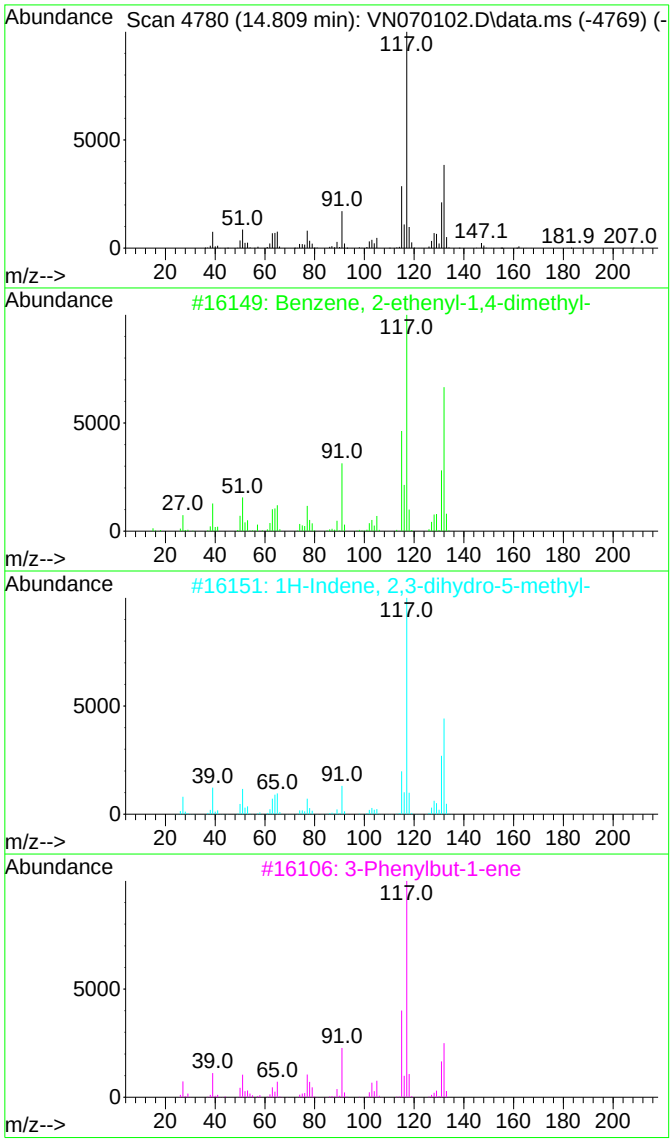
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

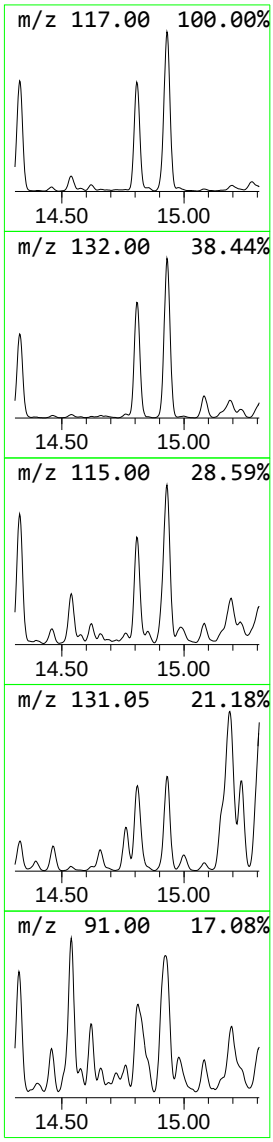
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.809	31.41 ug/l	5057180	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	94
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

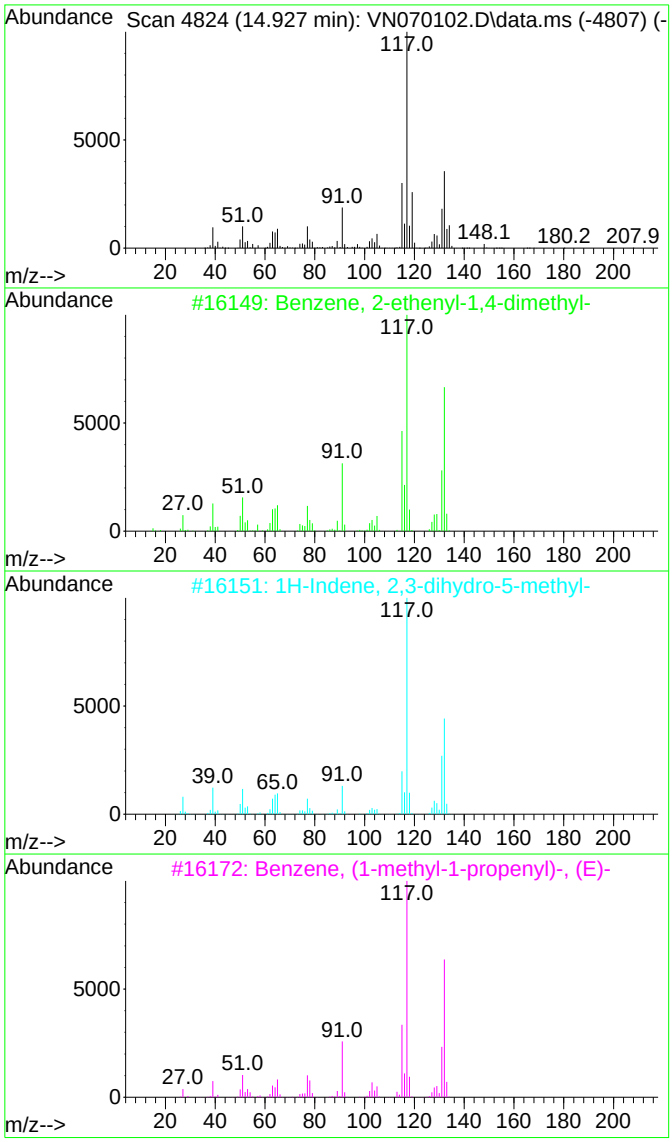
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

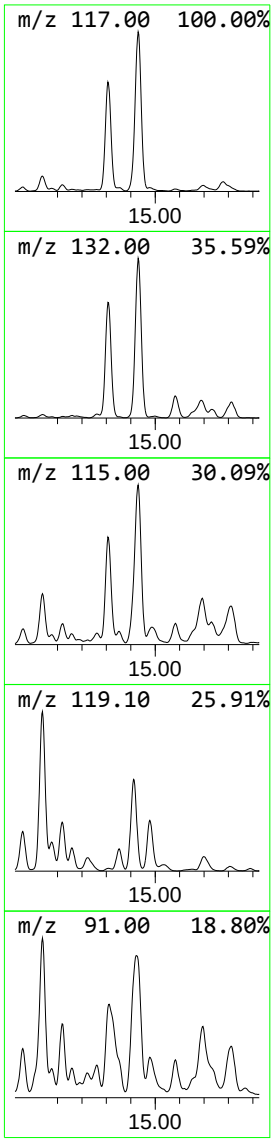
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	75.20 ug/l	12107800	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070102.D
Acq On : 15 Dec 2021 5:01
Operator : JC/MD
Sample : M5011-17 10X
Misc : 4.32g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 43 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-8-(10-10.5)

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.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, 2,2,3,3...	8.614	54.0	ug/l	8008000	2	8.968	7422100	50.0
Pentane, 2,3,4-...	9.883	46.9	ug/l	6968530	2	8.968	7422100	50.0
Pentane, 2,3,3-...	10.014	71.6	ug/l	10625200	2	8.968	7422100	50.0
Heptane, 3-methyl-	10.215	40.8	ug/l	6060920	2	8.968	7422100	50.0
Octane, 2-methyl-	11.527	37.5	ug/l	8036460	3	11.744	10709800	50.0
Benzene, 4-ethy...	14.214	81.2	ug/l	13077800	4	13.672	8050250	50.0
1-Phenyl-1-butene	14.324	37.8	ug/l	6082640	4	13.672	8050250	50.0
Benzene, 1,2,3,...	14.538	37.8	ug/l	6080650	4	13.672	8050250	50.0
Benzene, 2-ethe...	14.809	31.4	ug/l	5057180	4	13.672	8050250	50.0
1H-Indene, 2,3-...	14.927	75.2	ug/l	12107800	4	13.672	8050250	50.0