LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

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: VN070099.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.124	407	423	444	rBV3	179782	418121	1.29%	0.140%
2	5.074	1123	1150	1156	rBV4	164891	465848	1.43%	0.156%
3	5.616	1320	1352	1389	rBV4	233012	806825	2.48%	0.270%
4	6.114	1511	1538	1568	rBV4	224538	708278	2.18%	0.237%
5	6.898	1806	1830	1856	rBV7	115035	359085	1.11%	0.120%
6	7.042	1861	1884	1905	rBV2	591822	1779254	5.48%	0.595%
7	7.144	1906	1922	1941	rVB3	224662	612043	1.88%	0.205%
8	8.021	2228	2249	2257	rBV	901779	2151304	6.62%	0.720%
9	8.083	2260	2272	2289	rVV2	2083067	5379325	16.56%	1.800%
10	8.174	2290	2306	2330	rVB	1675868	4266850	13.13%	1.428%
11	8.295	2331	2351	2369	rBV2	768901	1761616	5.42%	0.589%
12	8.437	2385	2404	2435	rBV	1124128	2648482	8.15%	0.886%
13	8.614	2436	2470	2494	rBV	3009660	8591281	26.45%	2.874%
14	8.708	2496	2505	2526	rVB5	218699	572385	1.76%	0.191%
15	8.826	2527	2549	2579	rVB3	698083	1881060	5.79%	0.629%
16	8.965	2581	2601	2617	rBV2	3462490	7658795	23.58%	2.562%
17	9.124	2648	2660	2674	rVB2	222236	420279	1.29%	0.141%
18	9.201	2674	2689	2696	rBV5	177480	382073	1.18%	0.128%
19	9.456	2751	2784	2796	rBV2	1575265	3948851	12.16%	1.321%
20	9.510	2796	2804	2821	rVB2	761295	1471954	4.53%	0.492%
21	9.593	2821	2835	2843	rBV2	222121	450543	1.39%	0.151%
22	9.639	2845	2852	2866	rVB3	243117	445287	1.37%	0.149%
23	9.735	2867	2888	2902	rBV5	213256	537310	1.65%	0.180%
24	9.883	2927	2943	2964	rBV	2522981	5137264	15.81%	1.719%
25	9.974	2965	2977	2979	rBV2	662871	881717	2.71%	0.295%
26	10.014	2981	2992	3006	rVB	3268009	6409941	19.73%	2.145%
27	10.078	3007	3016	3025	rBV2	621034	993525	3.06%	0.332%
28	10.215	3047	3067	3092	rBV2	824584	2118684	6.52%	0.709%
29	10.320	3093	3106	3112	rBV4	244535	455279	1.40%	0.152%
30	10.371	3112	3125	3137	rVB	2053163	3631593	11.18%	1.215%
31	10.440	3137	3151	3164	rBV	5009756	9217886	28.38%	3.084%
32	10.505	3166	3175	3189	rVB	990400	1778205	5.47%	0.595%
33	10.625	3204	3220	3233	rVB4	1142763	2430196	7.48%	0.813%
34	10.861	3295	3308	3313	rBV2	758602	1382406	4.26%	0.463%
35	10.963	3336	3346	3358	rVB5	257392	495270	1.52%	0.166%
36	11.052	3359	3379	3391	rBV3	275850	635070	1.95%	0.212%

LSC Area Percent Report

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Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
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Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

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	11.154	3393	3417	3432	rBV4	447123	1402827	4.32%	0.469%
38	11.253	3447	3454	3464	rVB4	263928	375148	1.15%	0.126%
39	11.454	3521	3529	3538	rVB2	237869	328183	1.01%	0.110%
40	11.527	3546	3556	3582	rVB5	881962	2192960	6.75%	0.734%
41	11.626	3582	3593	3615	rBV	701767	1403177	4.32%	0.469%
42	11.744	3619	3637	3652	rVV	5616754	10364813	31.91%	3.468%
43	11.843	3655	3674	3694	rVV	4372314	8207370	25.27%	2.746%
44	11.950	3696	3714	3740	rVB	13945679	28428480	87.51%	9.511%
45	12.184	3792	3801	3813	rVB4	336744	545601	1.68%	0.183%
46	12.278	3819	3836	3854	rVB	3272355	6236098	19.20%	2.086%
47	12.363	3862	3868	3879	rVB2	403195	627407	1.93%	0.210%
48	12.578	3937	3948	3958	rBV	1169314	2014605	6.20%	0.674%
49	12.728	3991	4004	4027	rVB2	4554368	10134636	31.20%	3.391%
50	12.873	4049	4058	4062	rBV3	351710	506970	1.56%	0.170%
51	12.919	4064	4075	4086	rVB	3849472	5938360	18.28%	1.987%
52	12.980	4086	4098	4106	rBV	12000247	19973881	61.49%	6.683%
53	13.058	4119	4127	4167	rVB	7149577	12798647	39.40%	4.282%
54	13.219	4172	4187	4204	rBV	4251740	7586009	23.35%	2.538%
55	13.366	4228	4242	4257	rBV	19826788	32484882	100.00%	10.868%
56	13.498	4275	4291	4295	rBV3	384285	876173	2.70%	0.293%
57	13.605	4325	4331	4338	rBV4	240846	326917	1.01%	0.109%
58	13.675	4345	4357	4362	rBV	4971821	8165395	25.14%	2.732%
59	13.838	4406	4418	4420	rBV2	1113390	1418107	4.37%	0.474%
60	13.871	4421	4430	4437	rVV3	4300352	7777529	23.94%	2.602%
61	13.906	4439	4443	4466	rVV4	3961919	7319611	22.53%	2.449%
62	13.991	4467	4475	4488	rVB3	744017	1131269	3.48%	0.378%
63	14.066	4494	4503	4514	rVB	856056	1335496	4.11%	0.447%
64	14.128	4515	4526	4532	rBV	1842065	2972995	9.15%	0.995%
65	14.214	4547	4558	4572	rVV	3196337	5068793	15.60%	1.696%
66	14.278	4575	4582	4588	rVV	349825	477861	1.47%	0.160%
67	14.324	4589	4599	4612	rVB3	1646185	2747214	8.46%	0.919%
68	14.458	4638	4649	4660	rBV	684947	1193036	3.67%	0.399%
69	14.538	4669	4679	4686	rBV	1539490	2404198	7.40%	0.804%
70	14.579	4686	4694	4703	rVV	2255899	3378984	10.40%	1.130%
71	14.622	4703	4710	4718	rVV2	881435	1255254	3.86%	0.420%
72	14.659	4719	4724	4739	rVB2	587916	851975	2.62%	0.285%
73	14.761	4753	4762	4770	rBV2	487584	712236	2.19%	0.238%
74	14.809	4770	4780	4788	rBV2	1290850	2188554	6.74%	0.732%
75	14.927	4807	4824	4836	rBV4	2046528	4861711	14.97%	1.627%
76	15.193	4902	4923	4934	rBV5	947897	2334127	7.19%	0.781%

LSC Area Percent Report

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HSB-5-(11.5-12)

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
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Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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

77	15.314	4954	4968	4982	rVB5	800003	1851185	5.70%	0.619%
78	15.504	5029	5039	5052	rVB	811939	1458214	4.49%	0.488%
79	15.611	5068	5079	5090	rBV3	293923	547431	1.69%	0.183%
80	15.804	5135	5151	5168	rBV2	373985	820360	2.53%	0.274%
81	16.617	5439	5454	5485	rVB	231939	588301	1.81%	0.197%

Sum of corrected areas: 298896865

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Data File : VN070099.D
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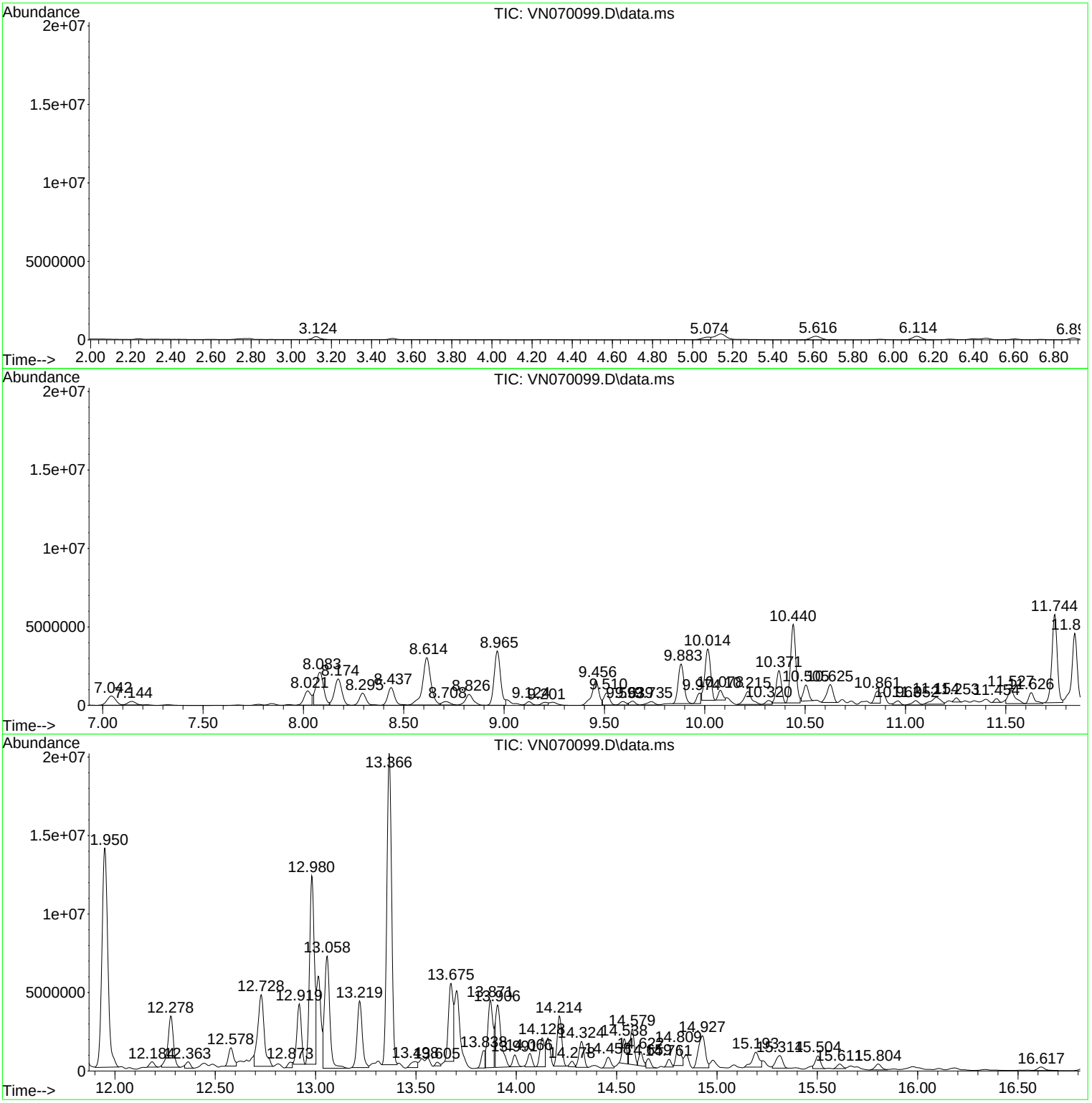
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

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 : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
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ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

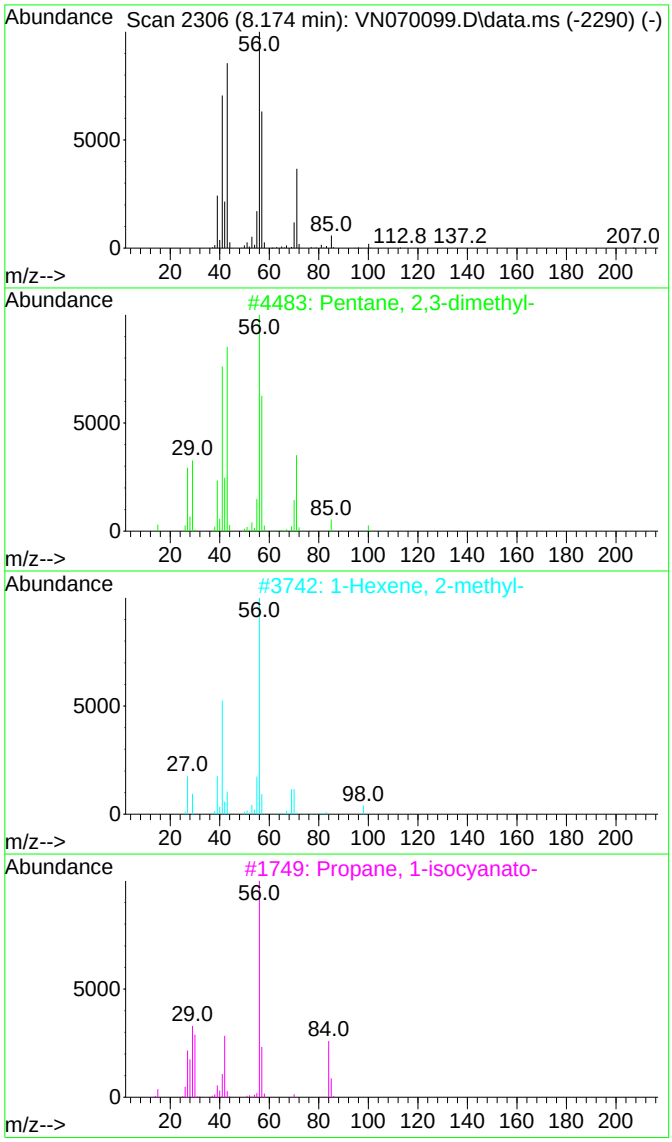
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

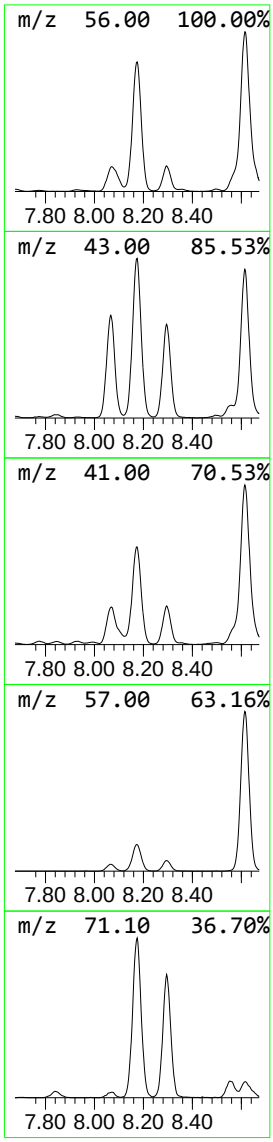
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Pentane, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.174	39.66 ug/l	4266850	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	1-Hexene, 2-methyl-	98	C7H14	006094-02-6	47
3	Propane, 1-isocyanato-	85	C4H7NO	000110-78-1	40
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	37
5	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

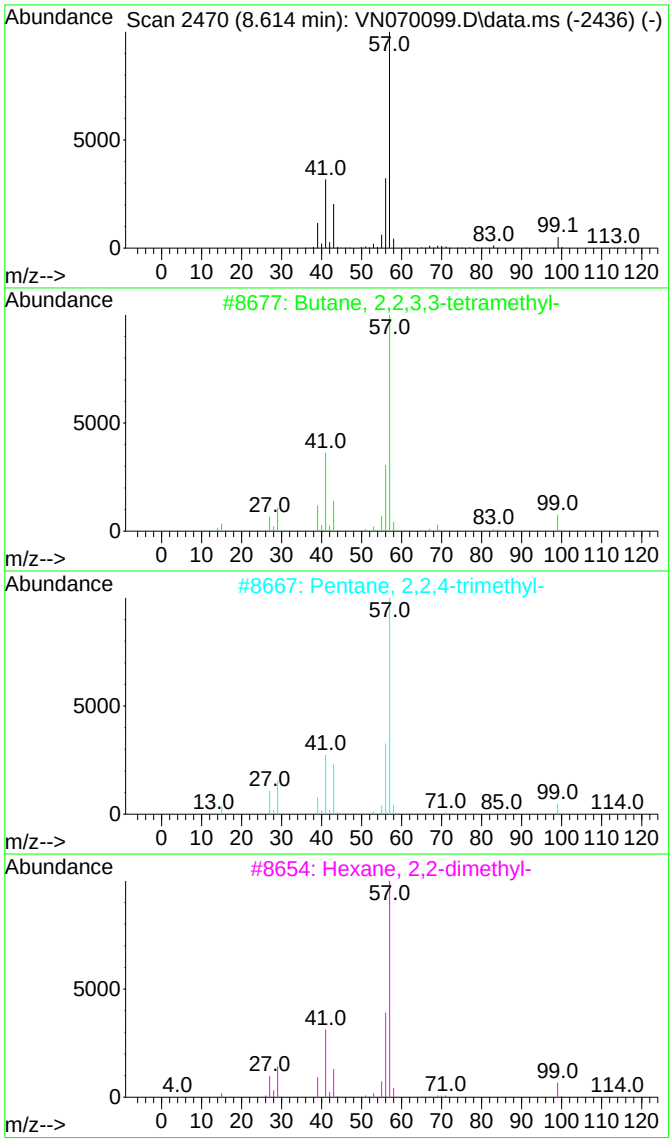
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

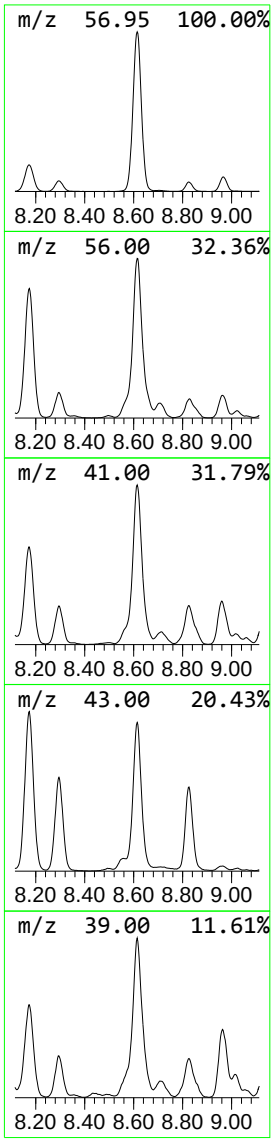
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.614	56.09 ug/l	8591280	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	72
2	Pentane, 2,2,4-trimethyl-	114	C8H18	000540-84-1	72
3	Hexane, 2,2-dimethyl-	114	C8H18	000590-73-8	64
4	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	59
5	Pentane, 2,2,4,4-tetramethyl-	128	C9H20	001070-87-7	59





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ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

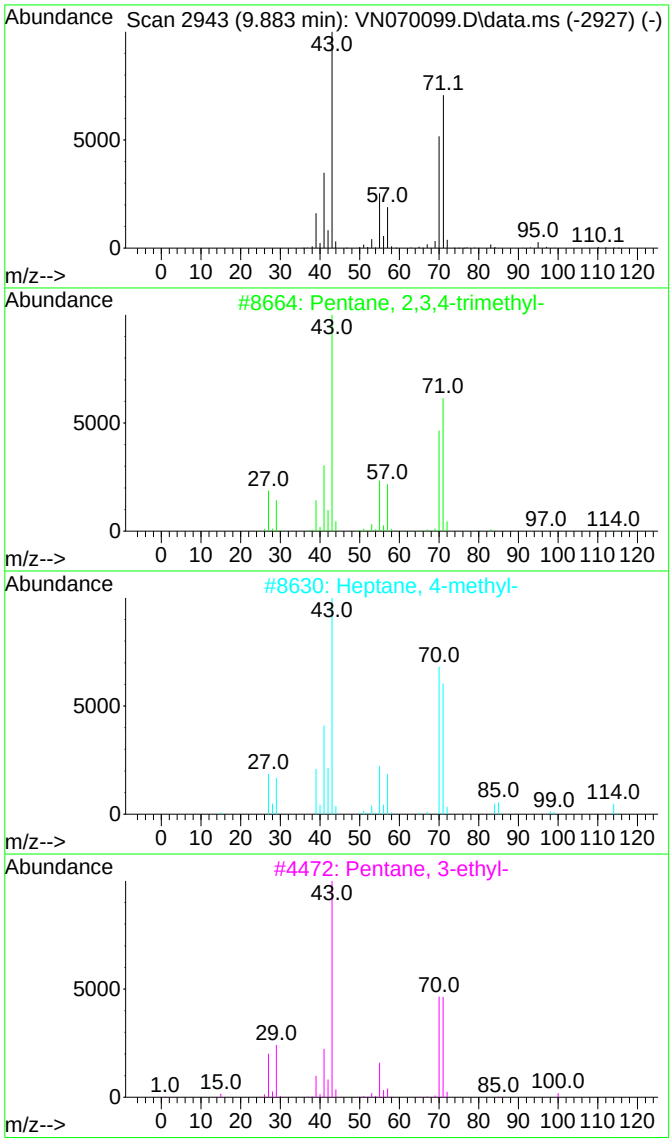
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

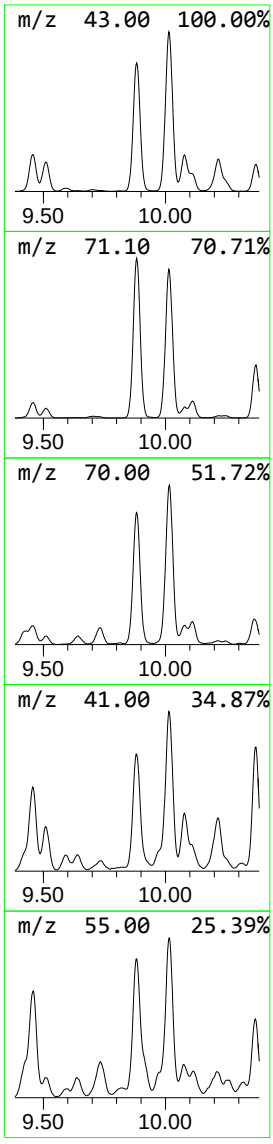
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane, 2,3,4-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.883	33.54 ug/l	5137260	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	78
4	Hexane, 2,3-dimethyl-	114	C8H18	000584-94-1	74
5	Hexane, 3,3,4-trimethyl-	128	C9H20	016747-31-2	56





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Operator : JC/MD
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Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

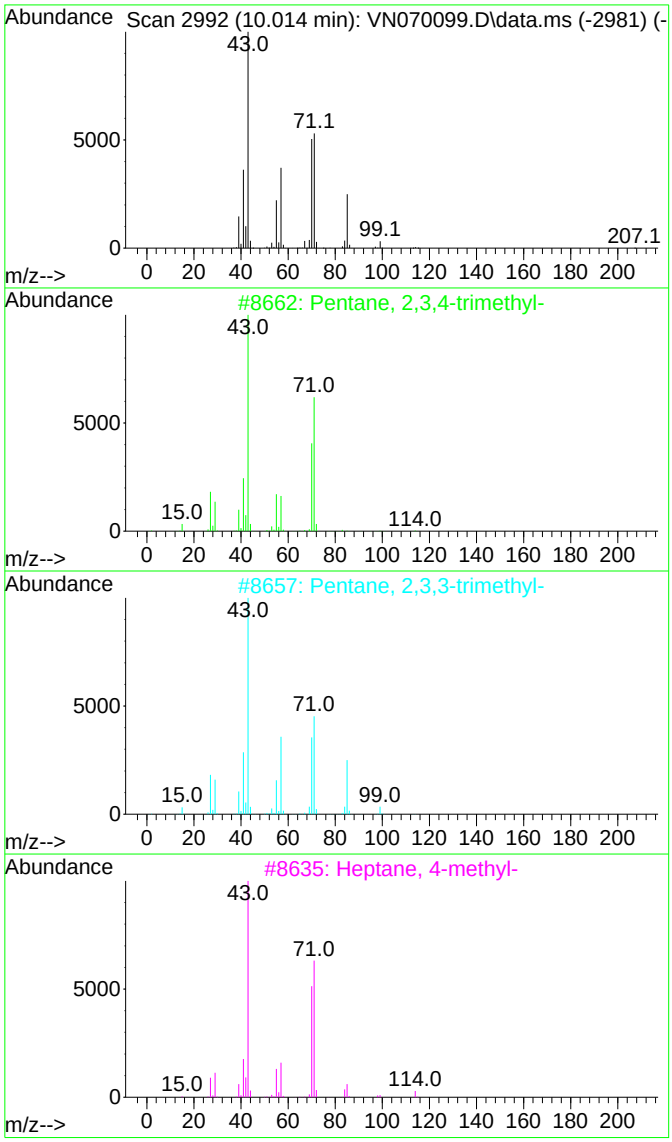
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

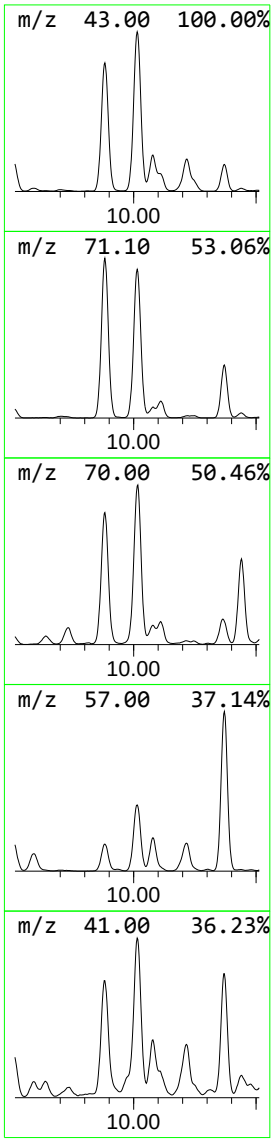
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.014	41.85 ug/l	6409940	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	72
2	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	72
3	Heptane, 4-methyl-	114	C8H18	000589-53-7	64
4	Decane, 3,3,4-trimethyl-	184	C13H28	049622-18-6	64
5	Pentane, 3-ethyl-	100	C7H16	000617-78-7	59





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Operator : JC/MD
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ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

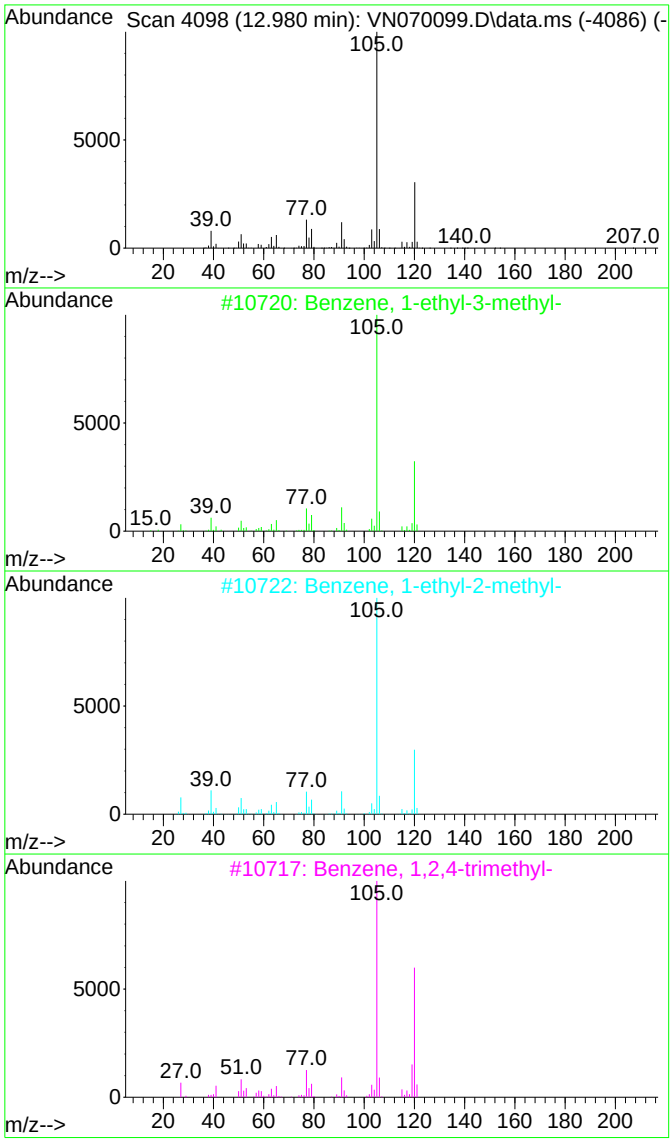
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

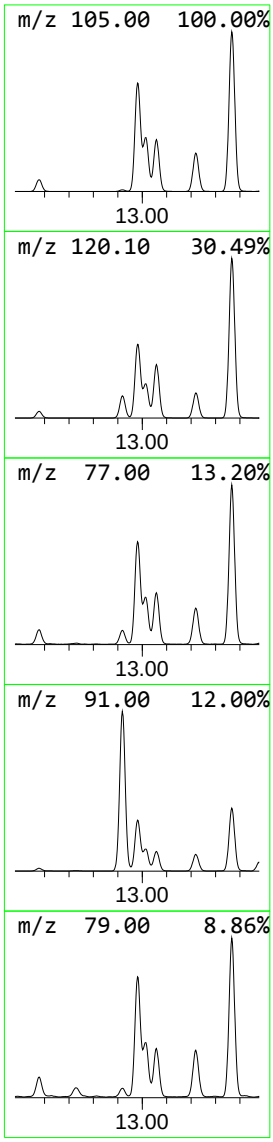
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1-ethyl-3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	122.31 ug/l	19973900	1,4-Dichlorobenzene-d4	13.675

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, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
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Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

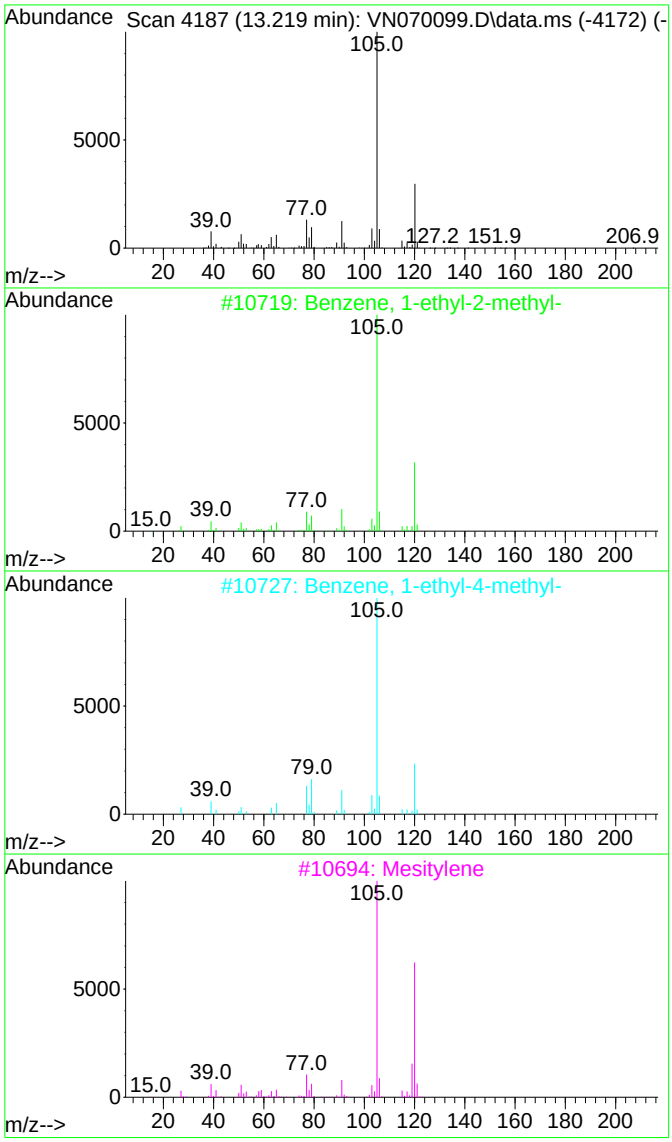
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

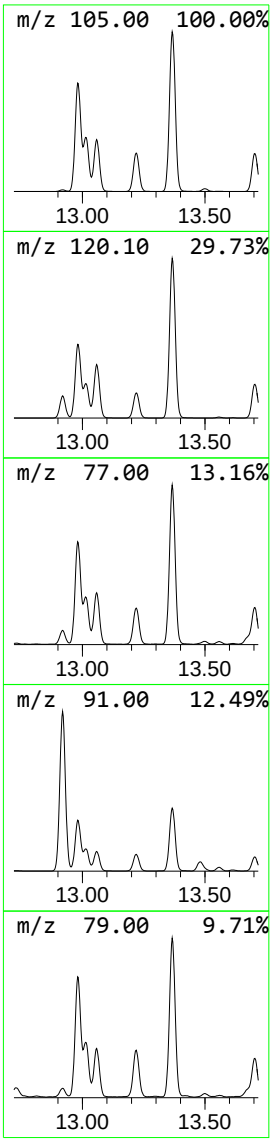
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.219	46.45 ug/l	7586010	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

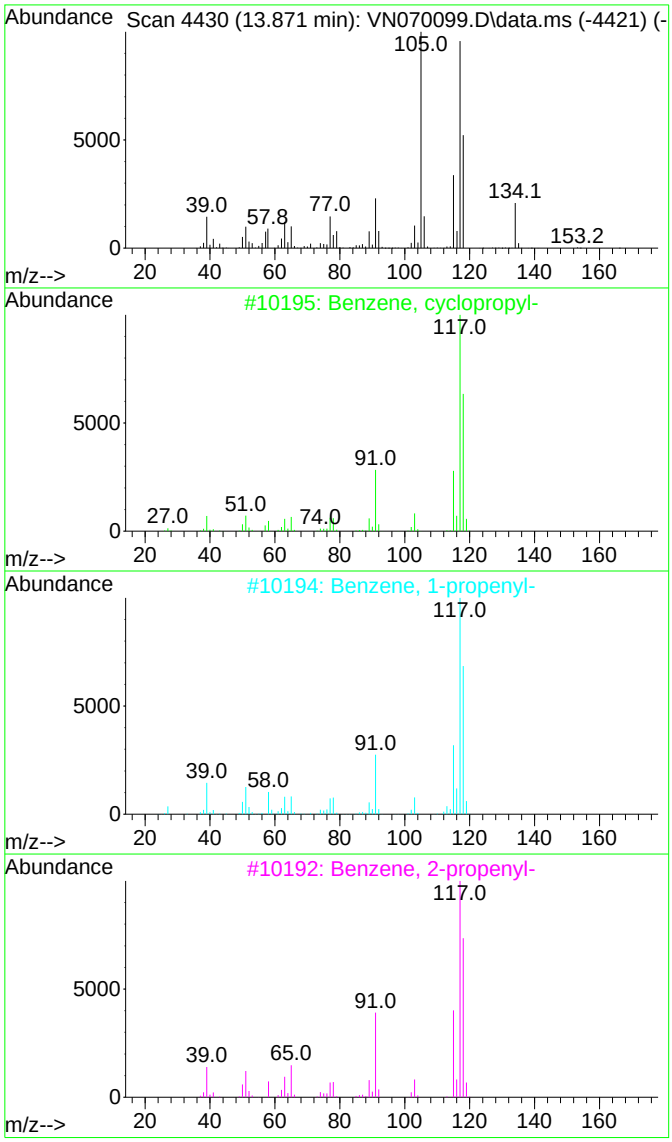
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

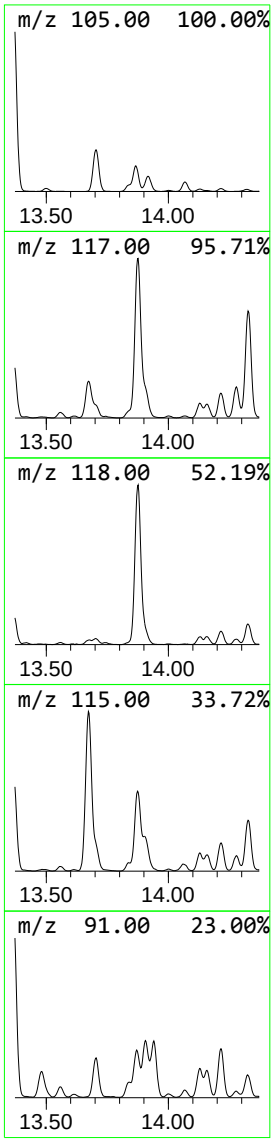
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, cyclopropyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.871	47.62 ug/l	7777530	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
2	Benzene, 1-propenyl-	118	C9H10	000637-50-3	64
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	64
4	Indane	118	C9H10	000496-11-7	60
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

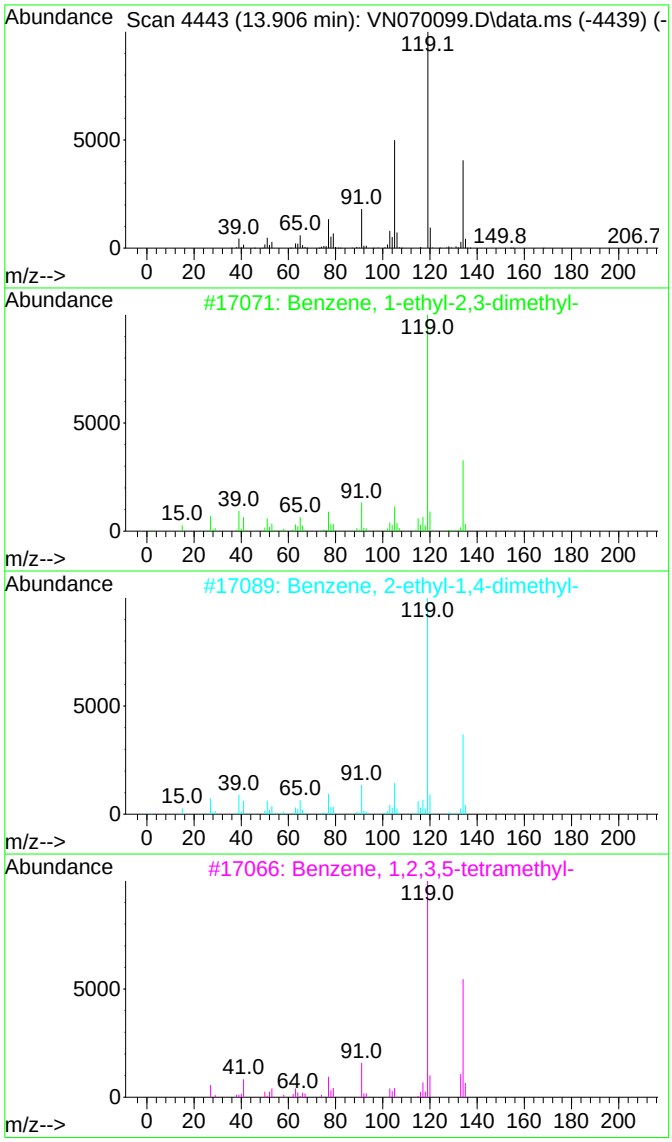
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

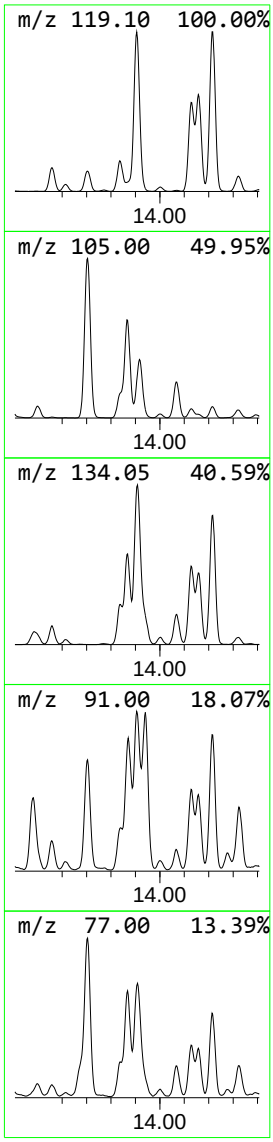
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.906	44.82 ug/l	7319610	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	90
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

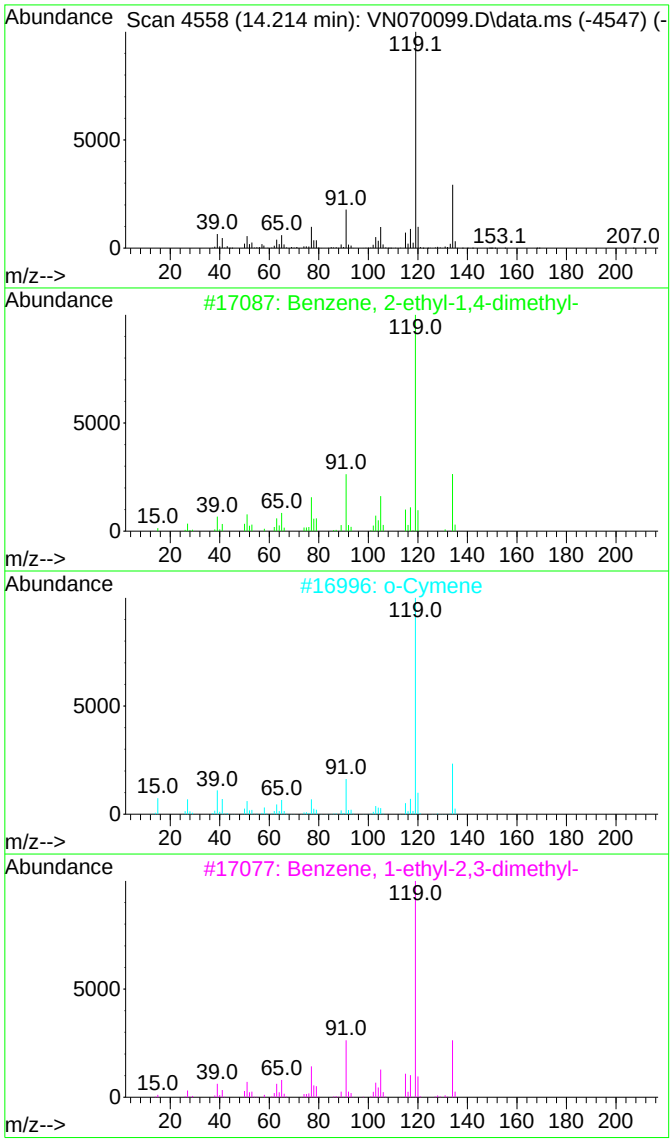
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

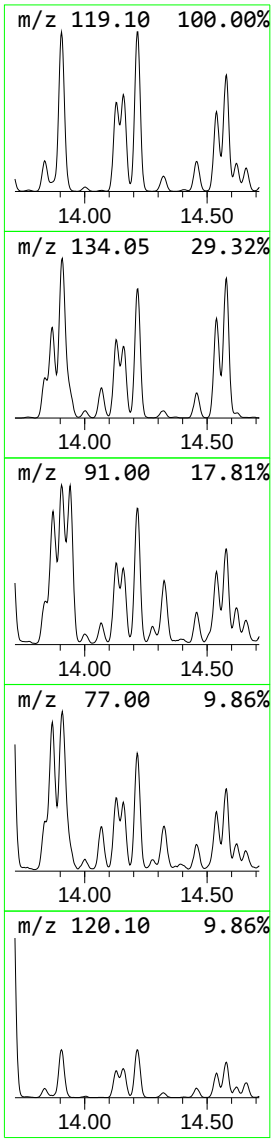
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.214	31.04 ug/l	5068790	1,4-Dichlorobenzene-d4	13.675

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

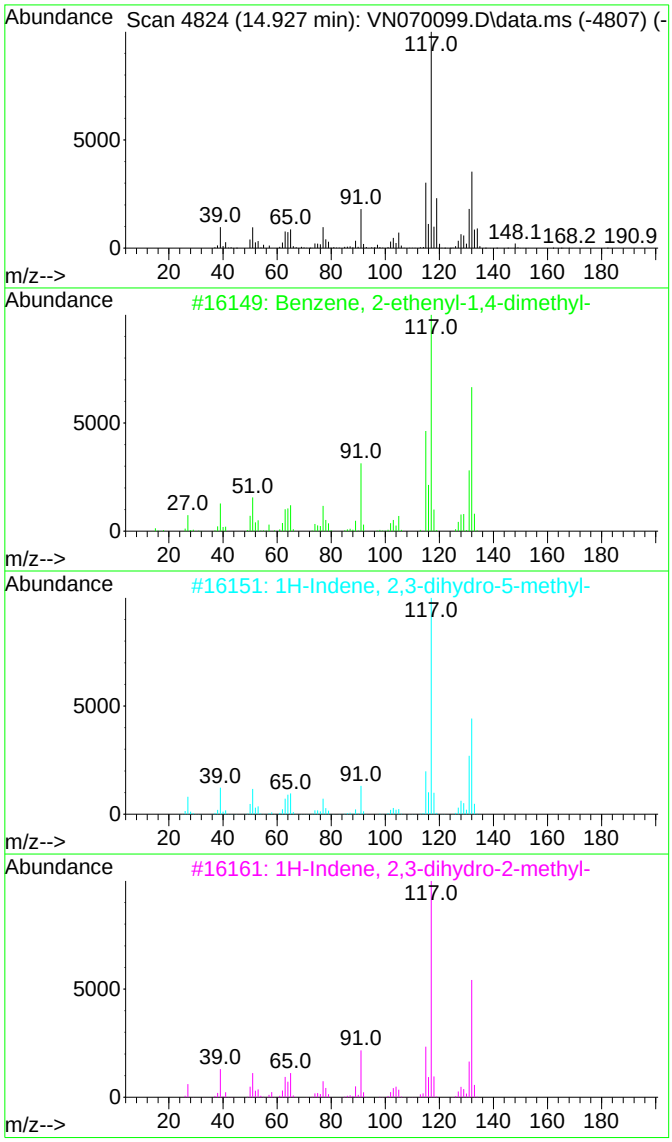
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N120221W.M
Quant Title : SW846 8260

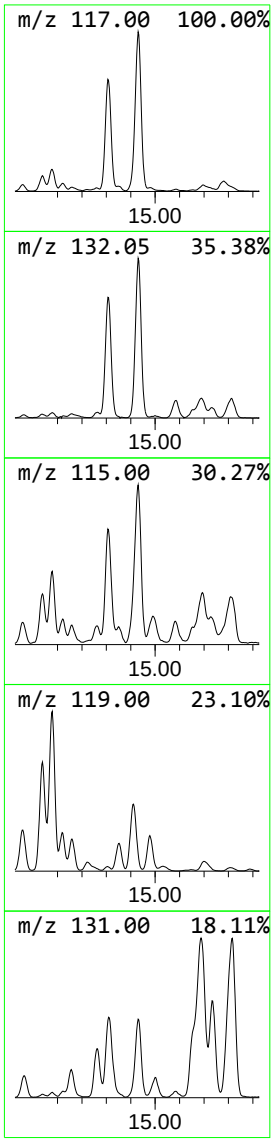
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	29.77 ug/l	4861710	1,4-Dichlorobenzene-d4	13.675

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	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	83
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	81
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN121421\
Data File : VN070099.D
Acq On : 15 Dec 2021 3:44
Operator : JC/MD
Sample : M5011-11 10X
Misc : 3.90g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 40 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
HSB-5-(11.5-12)

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
Pentane, 2,3-di...	8.174	39.7	ug/l	4266850	1	8.086	5379330	50.0
Butane, 2,2,3,3...	8.614	56.1	ug/l	8591280	2	8.968	7658800	50.0
Pentane, 2,3,4-...	9.883	33.5	ug/l	5137260	2	8.968	7658800	50.0
Pentane, 2,3,3-...	10.014	41.9	ug/l	6409940	2	8.968	7658800	50.0
Benzene, 1-ethy...	12.980	122.3	ug/l	19973900	4	13.675	8165400	50.0
Benzene, 1-ethy...	13.219	46.5	ug/l	7586010	4	13.675	8165400	50.0
Benzene, cyclop...	13.871	47.6	ug/l	7777530	4	13.675	8165400	50.0
Benzene, 1-ethy...	13.906	44.8	ug/l	7319610	4	13.675	8165400	50.0
Benzene, 2-ethy...	14.214	31.0	ug/l	5068790	4	13.675	8165400	50.0
Benzene, 2-ethe...	14.927	29.8	ug/l	4861710	4	13.675	8165400	50.0