LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

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_U\Method\82U062722W.M
Title : SW846 8260

Signal : TIC: VU049595.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.073	514	539	551	rBV	269662	618286	18.01%	0.850%
2	3.353	608	626	645	rBV	196620	433214	12.62%	0.595%
3	3.691	714	731	751	rBV	168832	375766	10.95%	0.516%
4	3.970	801	818	837	rBV2	70349	171342	4.99%	0.235%
5	4.527	974	991	1008	rVV	346412	844662	24.60%	1.161%
6	5.289	1212	1228	1239	rBV	178812	378299	11.02%	0.520%
7	5.372	1239	1254	1271	rVV4	818563	2112842	61.54%	2.904%
8	5.459	1271	1281	1302	rVB2	136387	315975	9.20%	0.434%
9	5.588	1303	1321	1345	rBV	490708	1209375	35.23%	1.662%
10	5.700	1345	1356	1372	rVB	195572	387494	11.29%	0.533%
11	5.848	1387	1402	1412	rBV	287559	610368	17.78%	0.839%
12	5.919	1413	1424	1434	rVV2	226897	504816	14.70%	0.694%
13	5.986	1434	1445	1474	rVB2	400711	874745	25.48%	1.202%
14	6.247	1512	1526	1534	rBV	451511	858271	25.00%	1.180%
15	6.292	1535	1540	1549	rVB2	113099	196563	5.73%	0.270%
16	6.761	1662	1686	1709	rBV	1462663	3433119	100.00%	4.718%
17	6.864	1709	1718	1733	rVB2	99546	187343	5.46%	0.257%
18	6.980	1741	1754	1767	rBV	239917	443328	12.91%	0.609%
19	7.073	1768	1783	1797	rVB	233815	469782	13.68%	0.646%
20	7.234	1821	1833	1853	rVB3	272199	590913	17.21%	0.812%
21	7.398	1870	1884	1889	rBV	161801	306610	8.93%	0.421%
22	7.498	1905	1915	1922	rBV	360966	597459	17.40%	0.821%
23	7.536	1923	1927	1936	rVB	224571	308026	8.97%	0.423%
24	7.658	1956	1965	1974	rBV	338425	539639	15.72%	0.742%
25	7.758	1984	1996	2011	rVB2	402913	768071	22.37%	1.056%
26	7.854	2011	2026	2032	rBV2	726507	1360145	39.62%	1.869%
27	7.893	2033	2038	2051	rVB2	840754	1419227	41.34%	1.950%
28	8.025	2068	2079	2086	rBV6	193162	445108	12.97%	0.612%
29	8.096	2096	2101	2112	rVB	235070	368374	10.73%	0.506%
30	8.240	2131	2146	2161	rVB4	485135	1170059	34.08%	1.608%
31	8.369	2174	2186	2193	rBV2	245096	457988	13.34%	0.629%
32	8.411	2193	2199	2212	rVB3	136273	277204	8.07%	0.381%
33	8.536	2229	2238	2250	rVB3	120319	201023	5.86%	0.276%
34	8.636	2251	2269	2282	rBV	462248	882635	25.71%	1.213%
35	8.703	2282	2290	2296	rBV4	94156	150973	4.40%	0.207%
36	8.751	2296	2305	2312	rBV4	233590	444099	12.94%	0.610%

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Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

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_U\Method\82U062722W.M
Title : SW846 8260

37	8.803	2314	2321	2331	rVB7	251934	450906	13.13%	0.620%
38	8.864	2331	2340	2350	rBV	559158	886690	25.83%	1.219%
39	8.925	2350	2359	2369	rBV	624259	977958	28.49%	1.344%
40	9.070	2390	2404	2412	rBV4	263892	558216	16.26%	0.767%
41	9.121	2413	2420	2426	rVV	225713	306381	8.92%	0.421%
42	9.166	2426	2434	2441	rVV5	311313	530014	15.44%	0.728%
43	9.211	2441	2448	2457	rVB2	442706	707595	20.61%	0.972%
44	9.263	2457	2464	2472	rBV	201034	273022	7.95%	0.375%
45	9.333	2474	2486	2501	rVB	749424	1258451	36.66%	1.730%
46	9.417	2502	2512	2520	rBV	570014	921906	26.85%	1.267%
47	9.462	2520	2526	2536	rVB4	103059	185269	5.40%	0.255%
48	9.565	2552	2558	2569	rVB4	158176	265173	7.72%	0.364%
49	9.665	2572	2589	2595	rBV6	278530	775576	22.59%	1.066%
50	9.716	2600	2605	2613	rVV2	503797	826378	24.07%	1.136%
51	9.761	2613	2619	2638	rVB5	318227	721888	21.03%	0.992%
52	9.880	2645	2656	2670	rBV10	154315	419784	12.23%	0.577%
53	10.025	2689	2701	2713	rBV6	441113	1087397	31.67%	1.494%
54	10.121	2723	2731	2742	rBV6	184751	417337	12.16%	0.574%
55	10.253	2759	2772	2794	rBV4	894123	2274465	66.25%	3.126%
56	10.359	2795	2805	2812	rBV3	725388	1375845	40.08%	1.891%
57	10.481	2833	2843	2852	rVB3	243553	469417	13.67%	0.645%
58	10.558	2852	2867	2875	rBV5	325002	742093	21.62%	1.020%
59	10.626	2876	2888	2904	rVB4	903211	1950755	56.82%	2.681%
60	10.697	2904	2910	2916	rBV3	113869	164417	4.79%	0.226%
61	10.761	2922	2930	2938	rBV6	341115	639958	18.64%	0.880%
62	10.809	2939	2945	2962	rVB5	451621	869533	25.33%	1.195%
63	10.906	2968	2975	2981	rBV	212988	308624	8.99%	0.424%
64	11.063	3017	3024	3030	rBV3	419171	718330	20.92%	0.987%
65	11.189	3056	3063	3071	rBV5	97261	166582	4.85%	0.229%
66	11.247	3073	3081	3087	rBV5	147878	252835	7.36%	0.347%
67	11.314	3088	3102	3114	rVB8	339275	945082	27.53%	1.299%
68	11.375	3115	3121	3126	rBV3	169473	242683	7.07%	0.334%
69	11.417	3127	3134	3142	rVB	671506	924829	26.94%	1.271%
70	11.465	3143	3149	3154	rBV5	136499	218574	6.37%	0.300%
71	11.575	3176	3183	3188	rBV5	149968	245120	7.14%	0.337%
72	11.642	3198	3204	3215	rVB4	497385	812288	23.66%	1.116%
73	11.713	3218	3226	3232	rBV	568034	910368	26.52%	1.251%
74	11.812	3249	3257	3270	rVB3	634136	1333789	38.85%	1.833%
75	11.877	3271	3277	3295	rVB2	244970	559830	16.31%	0.769%
76	12.095	3337	3345	3351	rBV2	401658	652435	19.00%	0.897%

LSC Area Percent Report

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ALS Vial : 19 Sample Multiplier: 1

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ClientSampleId :
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Integration Parameters: RTEINT.P
Integrator: RTE
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Sampling : 1 Min Area: 3 % of largest Peak
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Title : SW846 8260

77	12.185	3365	3373	3390	rVB6	838886	1931716	56.27%	2.655%
78	12.378	3427	3433	3453	rVB3	475042	1033748	30.11%	1.421%
79	12.571	3488	3493	3500	rVB	434876	557383	16.24%	0.766%
80	12.812	3561	3568	3571	rBV6	253664	358870	10.45%	0.493%
81	12.931	3597	3605	3612	rBV3	526670	1026161	29.89%	1.410%
82	13.028	3629	3635	3651	rVB4	605093	1364356	39.74%	1.875%
83	13.111	3654	3661	3665	rBV4	211731	340520	9.92%	0.468%
84	13.452	3758	3767	3783	rBV3	1166171	2221428	64.71%	3.053%
85	13.587	3805	3809	3815	rBV2	195515	212117	6.18%	0.292%
86	13.838	3880	3887	3904	rBV6	295260	963128	28.05%	1.324%
87	14.121	3970	3975	3984	rVB7	314237	474587	13.82%	0.652%
88	15.021	4248	4255	4268	rBV5	301211	598765	17.44%	0.823%
89	15.221	4305	4317	4326	rBV2	656970	1266494	36.89%	1.741%
90	15.275	4328	4334	4343	rVB6	250472	366125	10.66%	0.503%
91	15.410	4368	4376	4388	rVB	765320	1215900	35.42%	1.671%
92	16.150	4598	4606	4613	rBV3	380844	651549	18.98%	0.895%
93	16.259	4633	4640	4650	rVB	694912	1094926	31.89%	1.505%
94	16.407	4678	4686	4692	rBV	787036	1224160	35.66%	1.682%
95	16.446	4693	4698	4715	rVB	732367	1200562	34.97%	1.650%
96	17.114	4901	4906	4920	rVB	196921	287916	8.39%	0.396%
97	17.327	4967	4972	4980	rVB2	240460	330892	9.64%	0.455%
98	17.375	4981	4987	5004	rVB	237492	410235	11.95%	0.564%
99	17.539	5031	5038	5048	rVB	232835	380340	11.08%	0.523%
100	17.606	5052	5059	5071	rVV3	108594	189911	5.53%	0.261%

Sum of corrected areas: 72762725

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Instrument :
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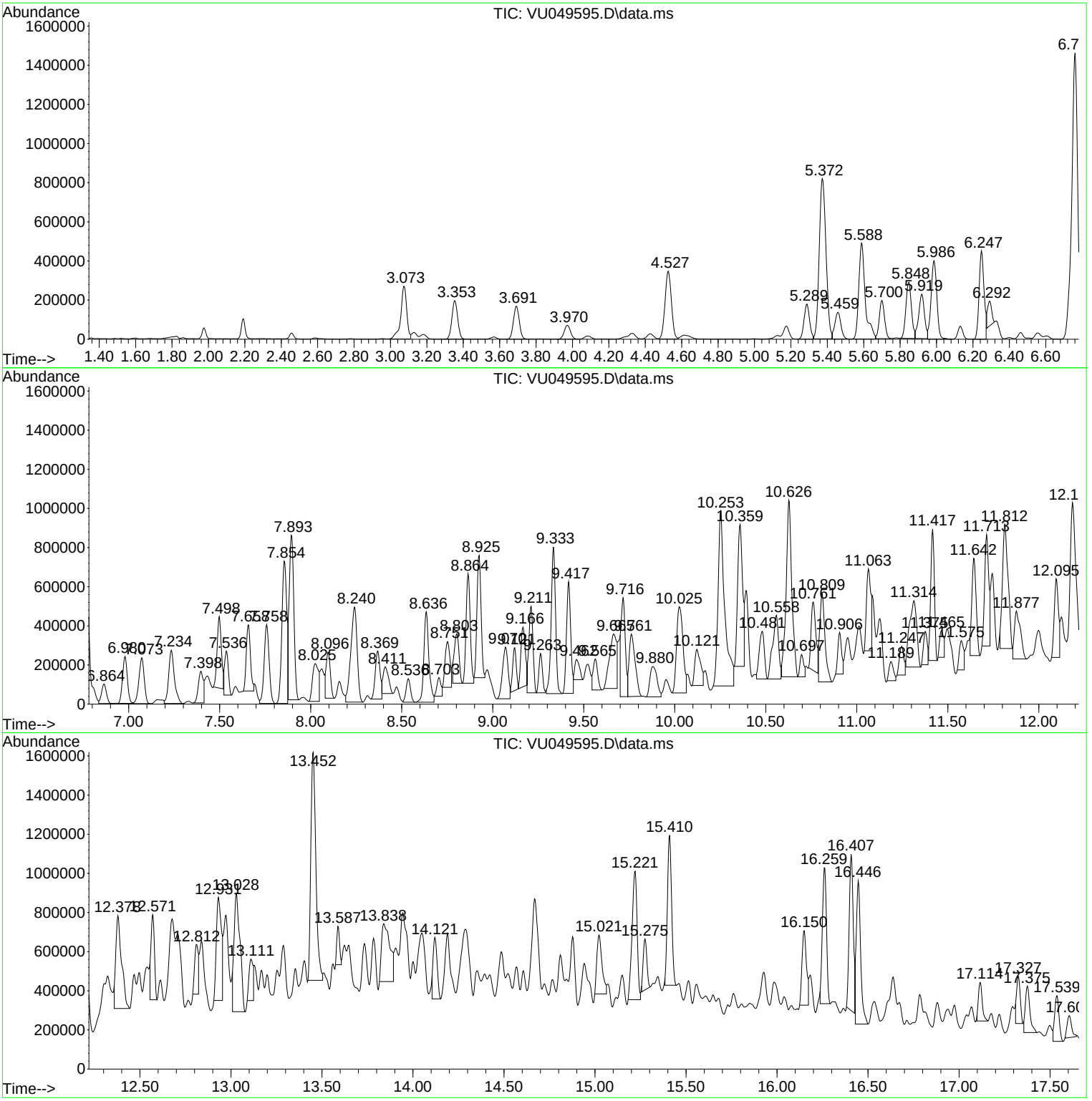
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
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Instrument :
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ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

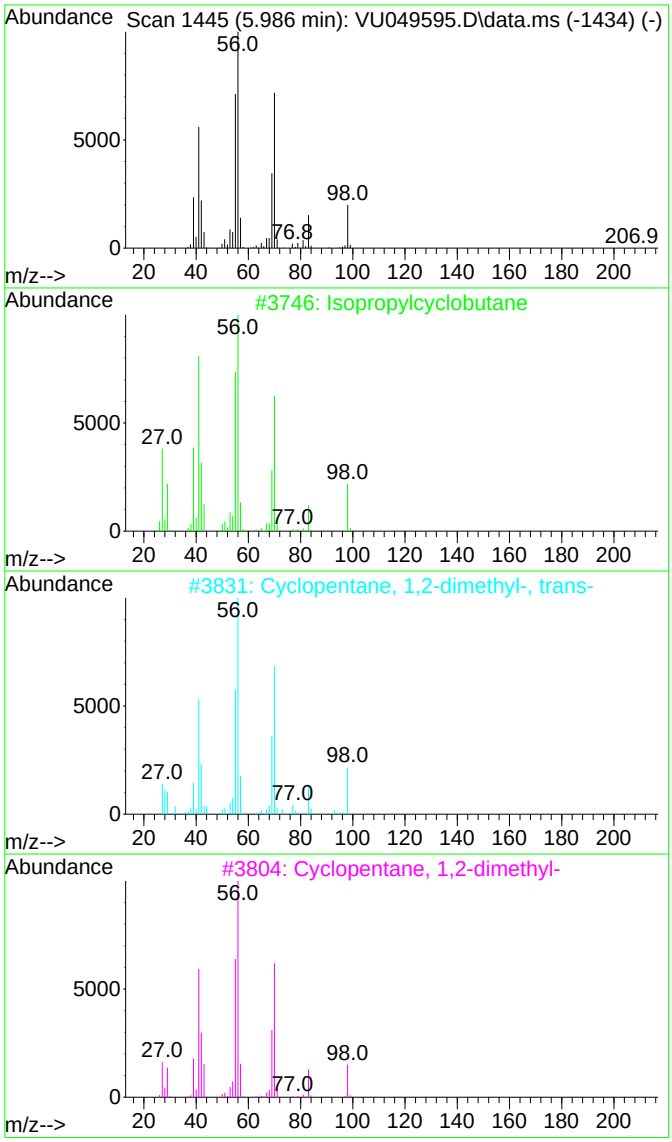
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

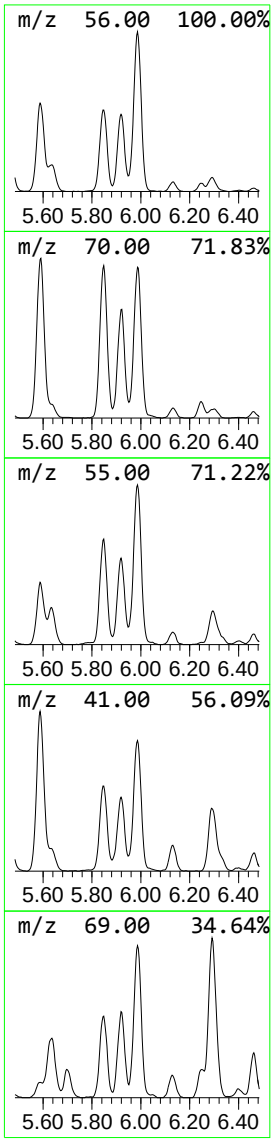
****

Peak Number 1 Isopropylcyclobutane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.986	50.96 ug/l	874745	1,4-Difluorobenzene	6.247

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-	98	C7H14	002452-99-5	87
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	78
5	1-Heptene	98	C7H14	000592-76-7	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
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Operator : SY/MD
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

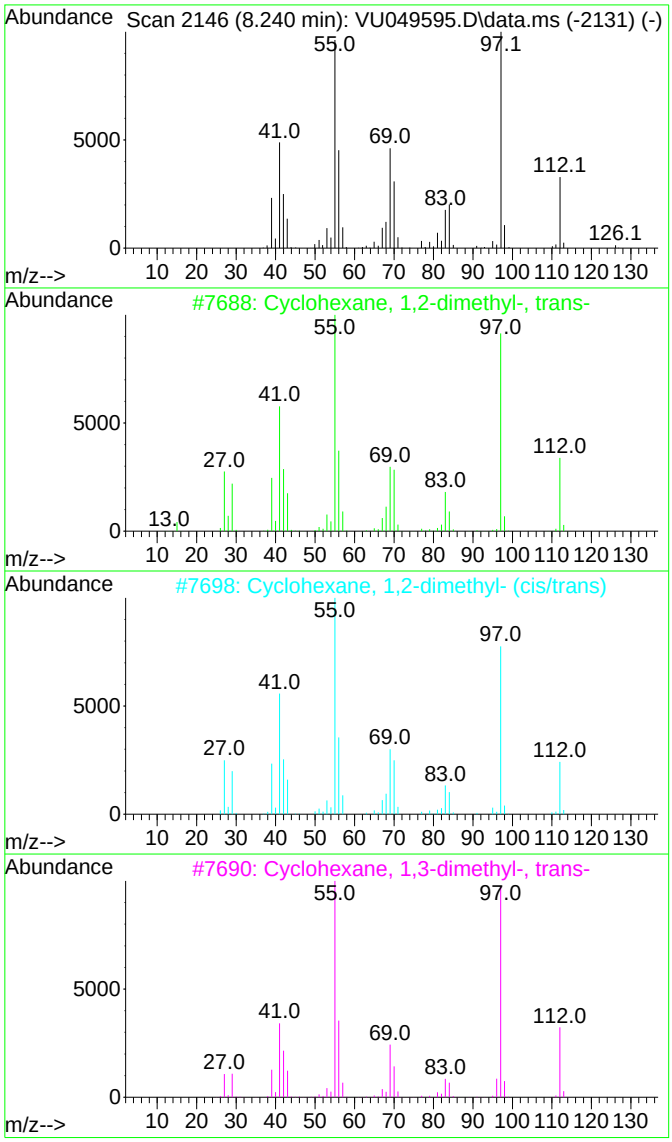
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

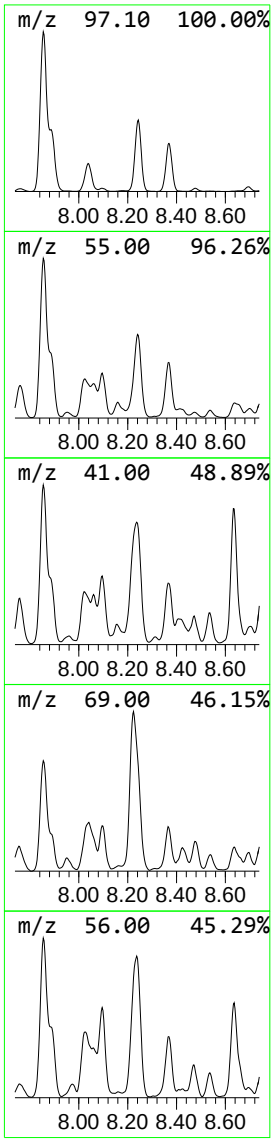
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclohexane, 1,2-dimethyl-,... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.240	63.46 ug/l	1170060	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	94
2	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	94
3	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	80
4	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	72
5	1H-Pyrazole, 4,5-dihydro-3,5,5-t...	112	C6H12N2	003975-85-7	70





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Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

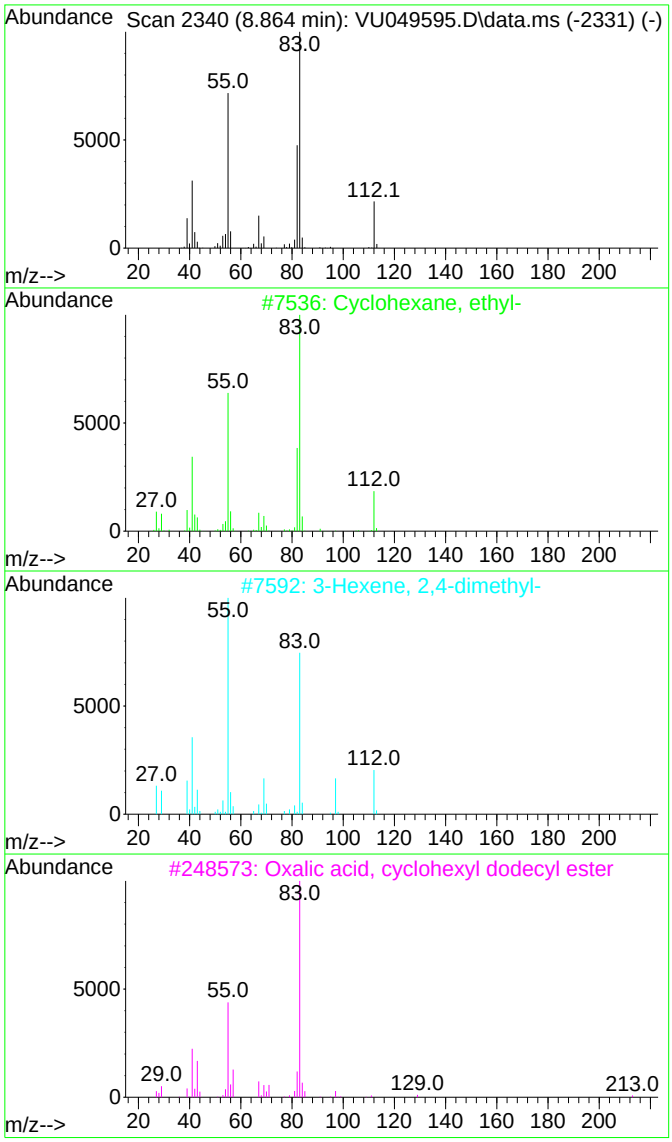
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

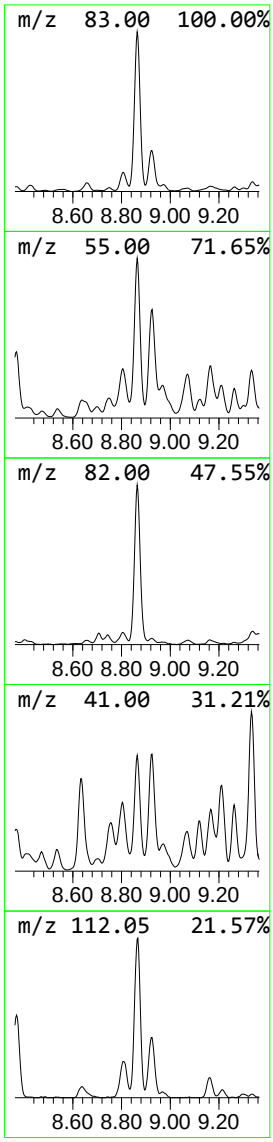
****

Peak Number 3 Cyclohexane, ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.864	48.09 ug/l	886690	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	86
2	3-Hexene, 2,4-dimethyl-	112	C8H16	082085-14-1	53
3	Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	53
4	Oxalic acid, dicyclohexyl ester	254	C14H22O4	1000309-30-9	50
5	Oxalic acid, cyclohexyl propyl e...	214	C11H18O4	1000309-30-3	50





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Operator : SY/MD
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

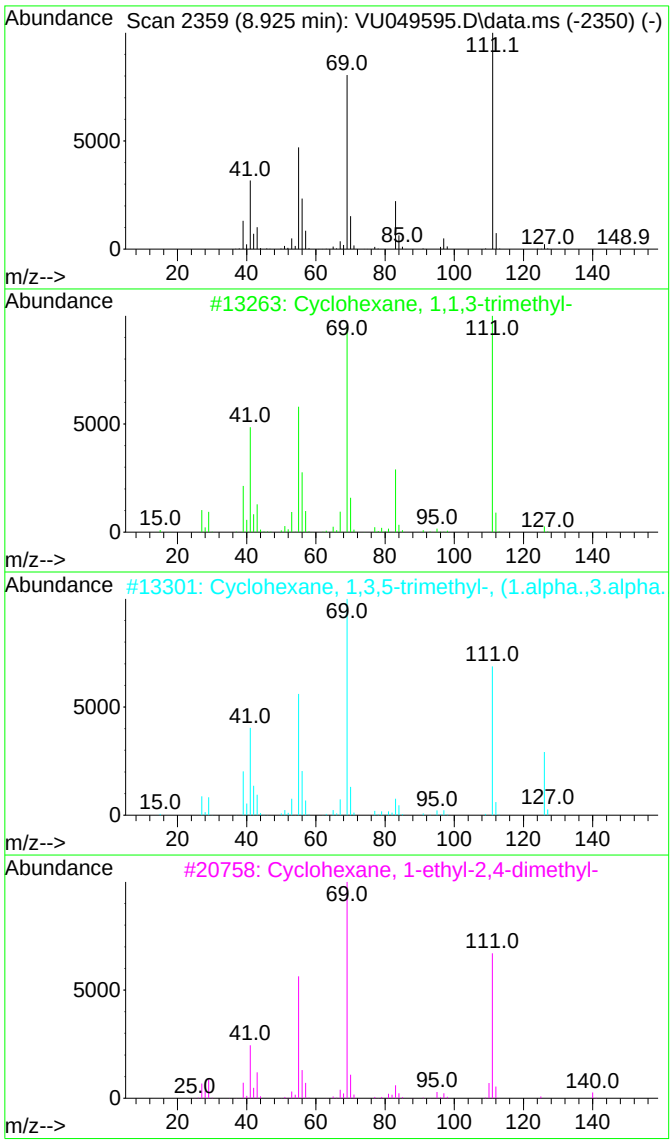
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

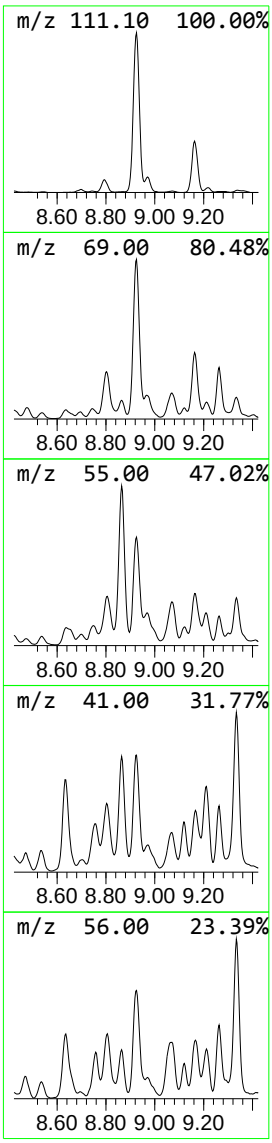
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Cyclohexane, 1,1,3-trimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.925	53.04 ug/l	977958	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	91
2	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-26-2	72
3	Cyclohexane, 1-ethyl-2,4-dimethyl-	140	C10H20	061142-69-6	72
4	Cyclohexane, 1,3,5-trimethyl-	126	C9H18	001839-63-0	56
5	Cyclohexane, 1,3,5-trimethyl-, (...	126	C9H18	001795-27-3	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
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ALS Vial : 19 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

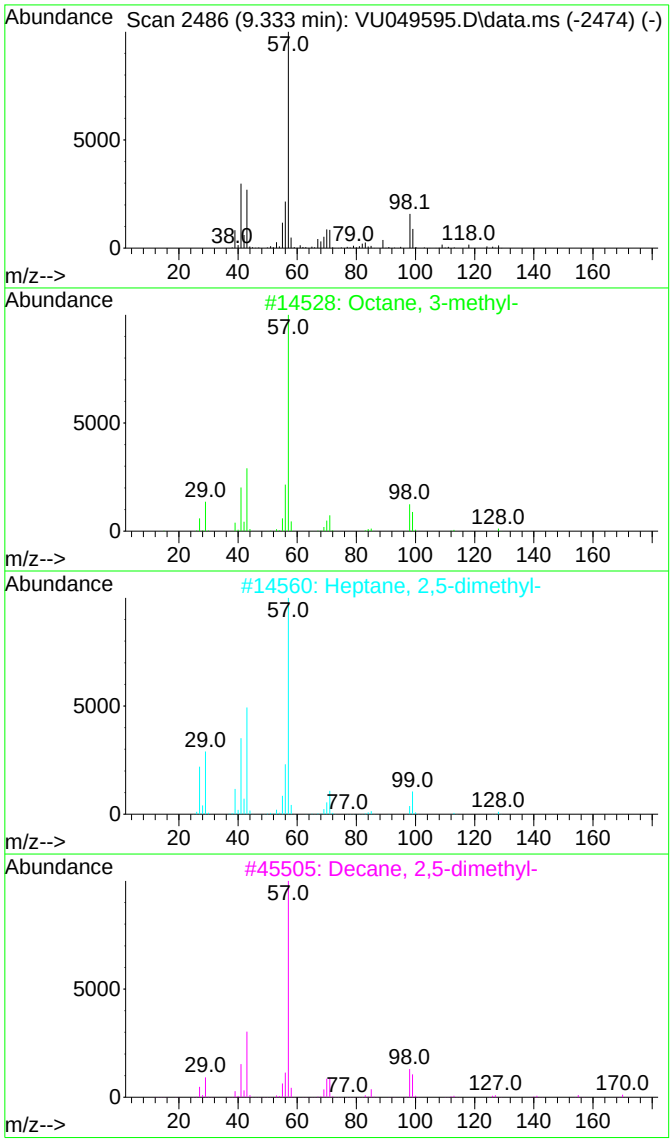
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

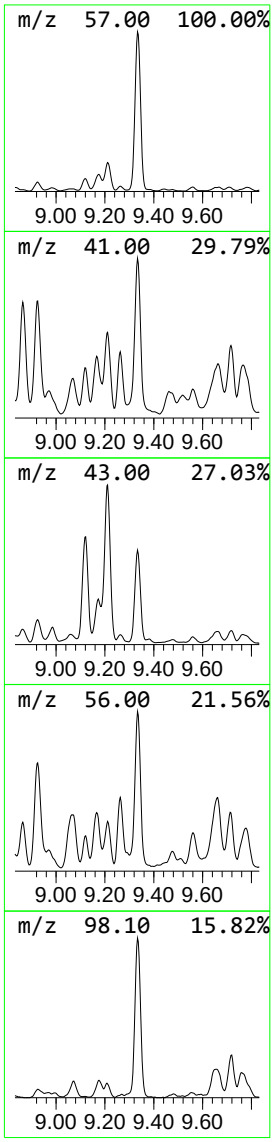
****

Peak Number 5 Octane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.333	68.25 ug/l	1258450	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-	128	C9H20	002216-33-3	83
2	Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	52
3	Decane, 2,5-dimethyl-	170	C12H26	017312-50-4	50
4	Undecane, 6-methyl-	170	C12H26	017302-33-9	50
5	Heptane, 4-propyl-	142	C10H22	003178-29-8	50





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Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

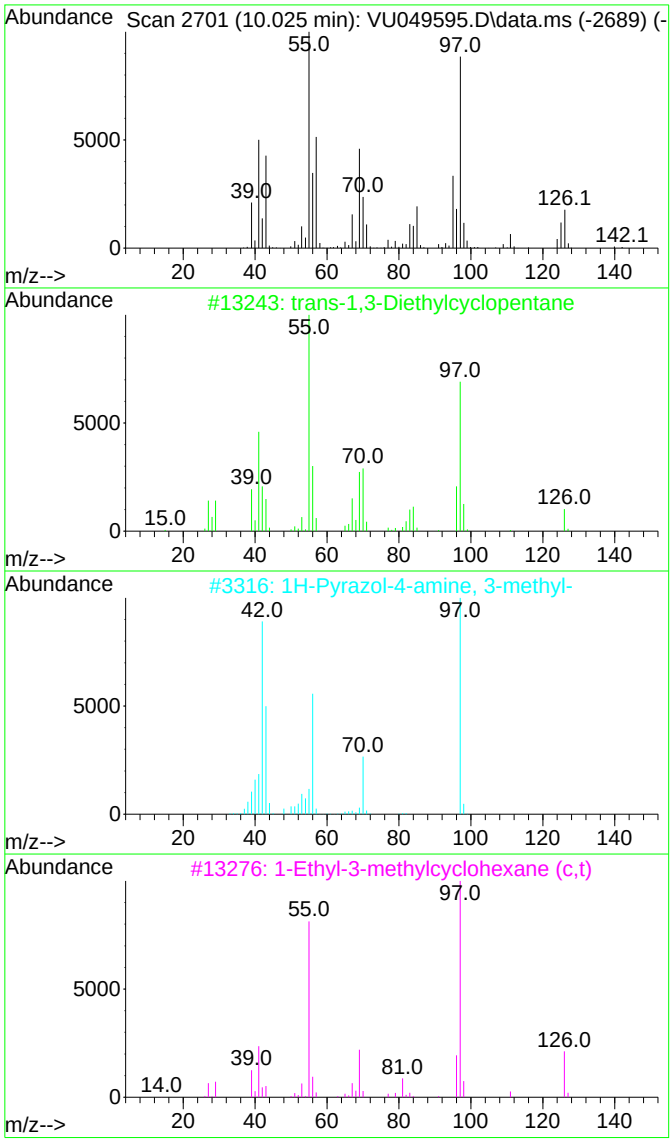
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

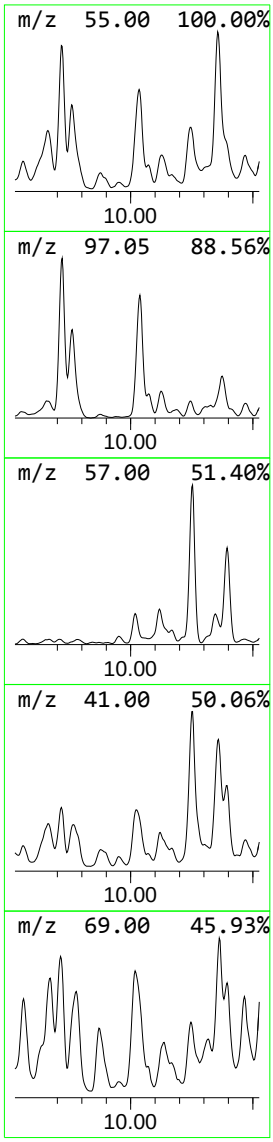
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 trans-1,3-Diethylcyclopentane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.025	58.98 ug/l	1087400	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	trans-1,3-Diethylcyclopentane	126	C9H18	1000113-87-1	76
2	1H-Pyrazol-4-amine, 3-methyl-	97	C4H7N3	1010338-28-2	49
3	1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	46
4	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	46
5	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	46





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Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
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Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

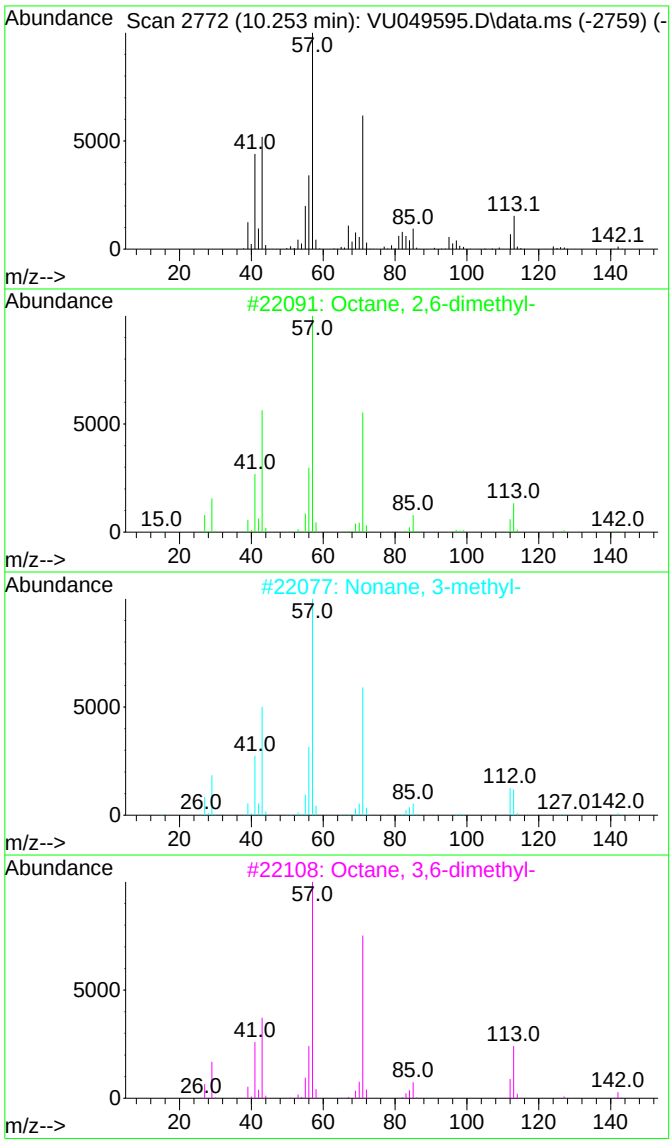
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

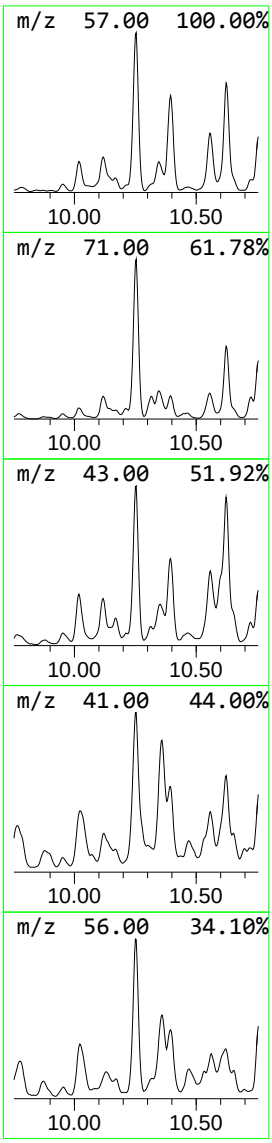
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Octane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.253	123.36 ug/l	2274470	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
2	Nonane, 3-methyl-	142	C10H22	005911-04-6	87
3	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	86
4	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	50
5	3-Bromooctane	192	C8H17Br	000999-64-4	47





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Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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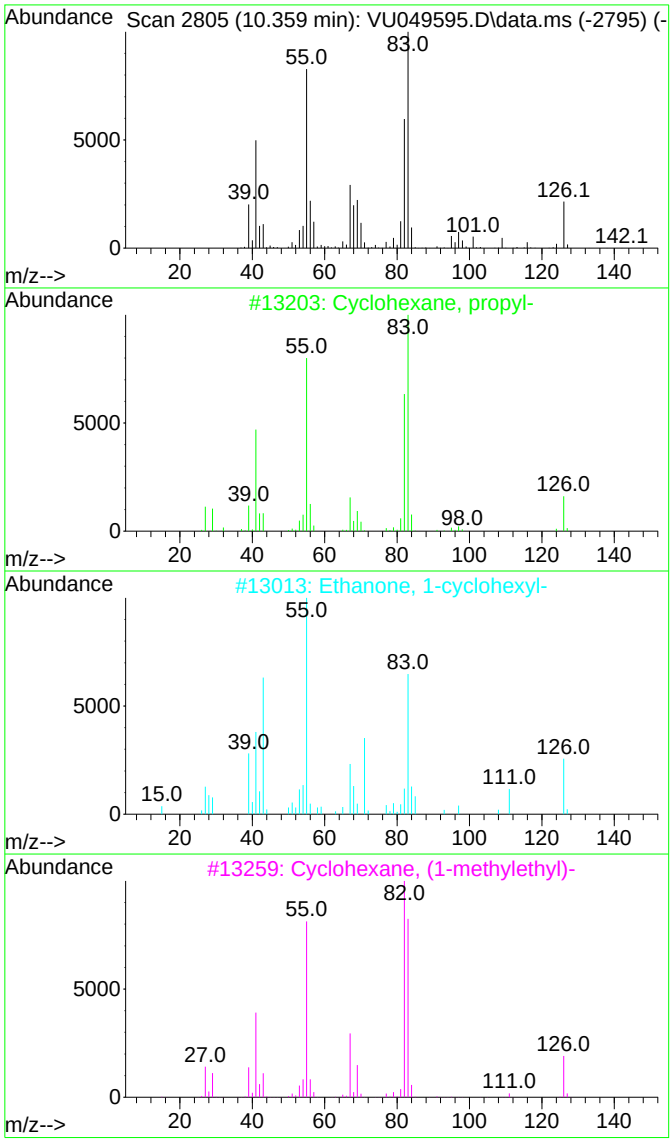
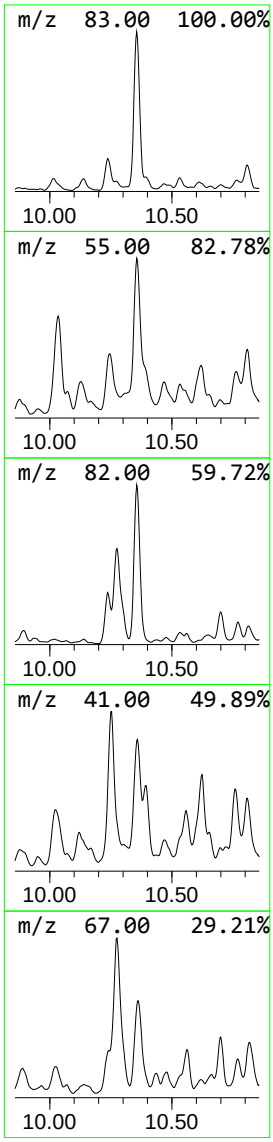
Peak Number 8 Cyclohexane, propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.359	74.62 ug/l	1375850	Chlorobenzene-d5	9.417

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	72
2	Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	58
3	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	58
4	Pyridine, 2,3,4,5-tetrahydro-	83	C5H9N	000505-18-0	50
5	Ethanone, 1-(1-methylcyclopentyl)-	126	C8H14O	013388-93-7	50





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Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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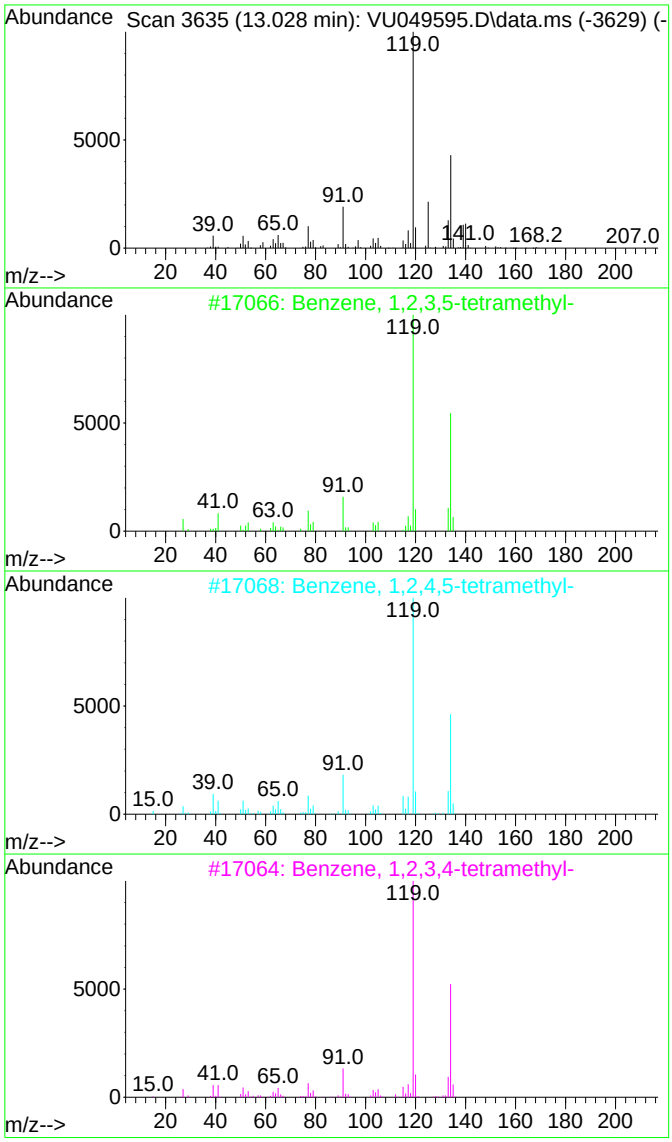
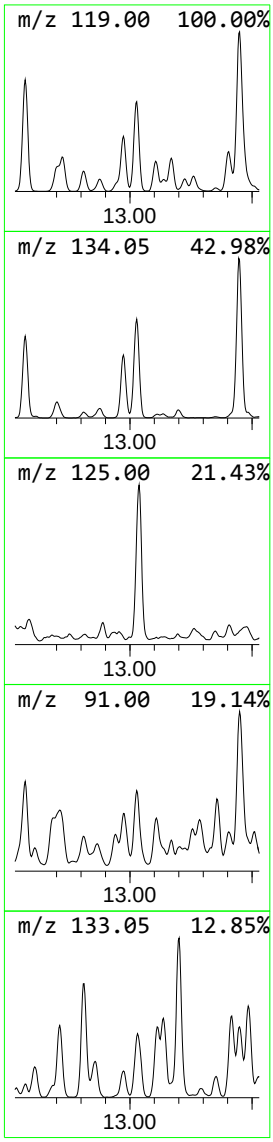
Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.028	51.15 ug/l	1364360	1,4-Dichlorobenzene-d4	11.813

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	94
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	p-Cymene	134	C10H14	000099-87-6	93





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Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.M
Quant Title : SW846 8260

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

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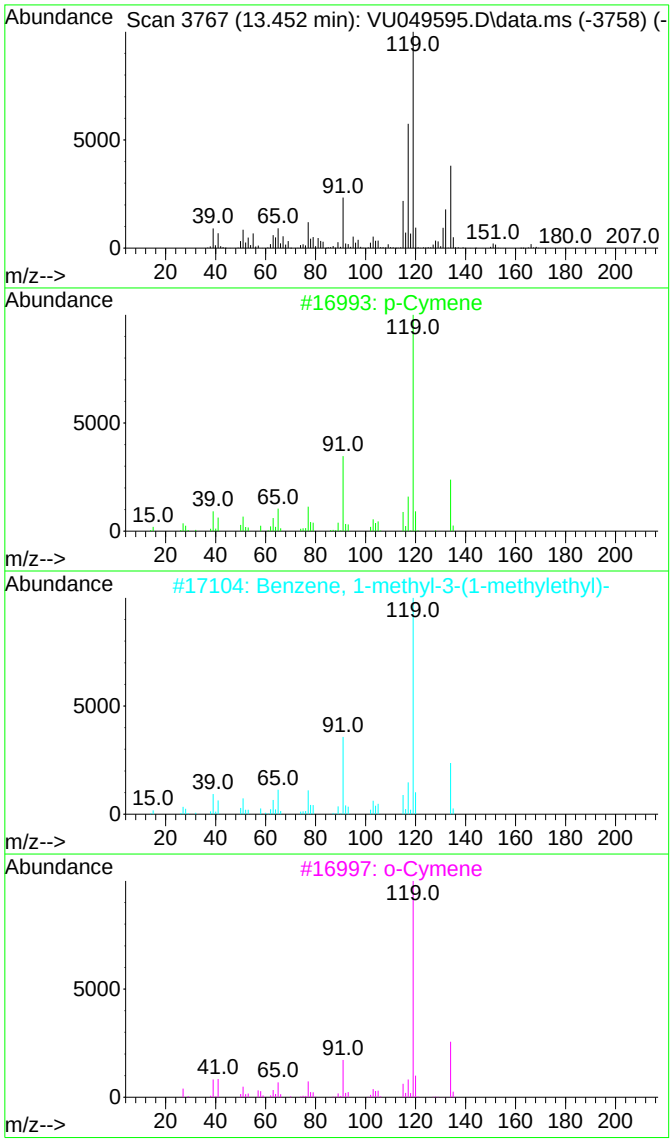
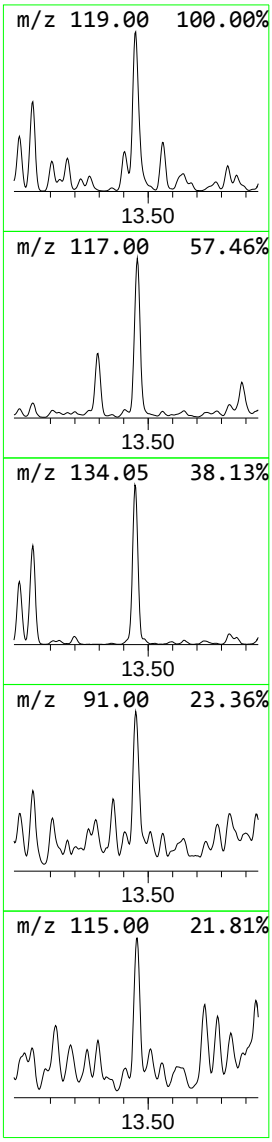
Peak Number 10 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.452	83.28 ug/l	2221430	1,4-Dichlorobenzene-d4	11.813

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	64
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64
3	o-Cymene	134	C10H14	000527-84-4	64
4	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	60
5	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	58





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU070122\
Data File : VU049595.D
Acq On : 01 Jul 2022 20:31
Operator : SY/MD
Sample : N3564-14 10X
Misc : 5.93g/5.0mL/100uL/5.0mL/MSVOA_U/MEOH
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Instrument :
MSVOA_U
ClientSampleId :
WC-14B-11-3-GR

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U062722W.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
Isopropylcyclob...	5.986	51.0	ug/l	874745	2	6.247	858271	50.0
Cyclohexane, 1,...	8.240	63.5	ug/l	1170060	3	9.417	921906	50.0
Cyclohexane, et...	8.864	48.1	ug/l	886690	3	9.417	921906	50.0
Cyclohexane, 1,...	8.925	53.0	ug/l	977958	3	9.417	921906	50.0
Octane, 3-methyl-	9.333	68.3	ug/l	1258450	3	9.417	921906	50.0
trans-1,3-Dieth...	10.025	59.0	ug/l	1087400	3	9.417	921906	50.0
Octane, 2,6-dim...	10.253	123.4	ug/l	2274470	3	9.417	921906	50.0
Cyclohexane, pr...	10.359	74.6	ug/l	1375850	3	9.417	921906	50.0
Benzene, 1,2,3,...	13.028	51.1	ug/l	1364360	4	11.813	1333790	50.0
o-Cymene	13.452	83.3	ug/l	2221430	4	11.813	1333790	50.0