LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
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_U\Method\82U092821W.M
Title : SW846 8260

Signal : TIC: VU044977.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.999	193	205	223	rBV	949209	1296797	6.36%	1.238%
2	2.208	259	270	292	rVB	1343912	1900247	9.31%	1.814%
3	2.321	295	305	319	rBV	157287	226674	1.11%	0.216%
4	3.089	519	544	554	rVV	875685	2161983	10.60%	2.064%
5	3.144	554	561	589	rVB2	297090	662869	3.25%	0.633%
6	3.369	609	631	654	rBV	580926	1297130	6.36%	1.238%
7	3.707	717	736	757	rBV	1048349	2329097	11.41%	2.223%
8	4.539	977	995	1017	rVB	1292896	3112959	15.26%	2.971%
9	5.295	1208	1230	1242	rBV	176425	410747	2.01%	0.392%
10	5.391	1242	1260	1274	rBV4	1614635	4240549	20.78%	4.048%
11	5.465	1276	1283	1307	rVB3	243251	597642	2.93%	0.570%
12	5.597	1309	1324	1333	rBV	516204	1138611	5.58%	1.087%
13	5.710	1349	1359	1366	rBV	173813	306889	1.50%	0.293%
14	5.780	1366	1381	1395	rVV	10192624	20405221	100.00%	19.478%
15	5.854	1395	1404	1415	rVV	578270	1252614	6.14%	1.196%
16	5.925	1416	1426	1436	rVV2	477746	1068123	5.23%	1.020%
17	5.996	1437	1448	1477	rVB	804171	1752004	8.59%	1.672%
18	6.137	1479	1492	1504	rBV	825648	1542238	7.56%	1.472%
19	6.253	1516	1528	1538	rBV	427283	806530	3.95%	0.770%
20	6.767	1662	1688	1710	rBV	3386711	7417515	36.35%	7.080%
21	6.986	1743	1756	1770	rVB	431632	780397	3.82%	0.745%
22	7.079	1770	1785	1799	rVB	225857	471299	2.31%	0.450%
23	7.240	1825	1835	1854	rVB	294796	621158	3.04%	0.593%
24	7.504	1905	1917	1925	rBV	254324	439940	2.16%	0.420%
25	7.767	1987	1999	2013	rVB6	92481	218522	1.07%	0.209%
26	7.861	2013	2028	2033	rBV2	726035	1295503	6.35%	1.237%
27	7.899	2034	2040	2052	rVV2	945636	1630518	7.99%	1.556%
28	7.973	2052	2063	2072	rVV	653221	1087195	5.33%	1.038%
29	8.038	2072	2083	2096	rVV5	203742	645448	3.16%	0.616%
30	8.102	2097	2103	2107	rVV	210263	335722	1.65%	0.320%
31	8.134	2108	2113	2125	rVB	310179	457001	2.24%	0.436%
32	8.247	2134	2148	2166	rVB2	508634	1037147	5.08%	0.990%
33	8.372	2171	2187	2196	rBV2	273992	521038	2.55%	0.497%
34	8.812	2313	2324	2332	rBV3	223171	395956	1.94%	0.378%
35	8.870	2332	2342	2351	rVB	594754	947030	4.64%	0.904%
36	8.928	2351	2360	2370	rBV	363769	592458	2.90%	0.566%

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ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
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_U\Method\82U092821W.M
Title : SW846 8260

37	9.337	2473	2487	2502	rVB2	142477	302980	1.48%	0.289%
38	9.420	2502	2513	2524	rBV	654920	1061911	5.20%	1.014%
39	9.513	2534	2542	2550	rVB	180219	280903	1.38%	0.268%
40	9.574	2550	2561	2574	rBV	1065329	1692705	8.30%	1.616%
41	9.697	2585	2599	2612	rBV	4644148	7755200	38.01%	7.403%
42	10.034	2689	2704	2712	rBV6	141087	342315	1.68%	0.327%
43	10.105	2720	2726	2753	rVB2	132799	323327	1.58%	0.309%
44	10.275	2754	2779	2791	rBV4	272573	694224	3.40%	0.663%
45	10.362	2796	2806	2826	rVB6	199749	475704	2.33%	0.454%
46	10.488	2834	2845	2854	rBV	313695	490492	2.40%	0.468%
47	10.635	2880	2891	2903	rVB	551102	930616	4.56%	0.888%
48	10.909	2966	2976	2986	rBV	395366	604126	2.96%	0.577%
49	11.002	2994	3005	3011	rBV	1211914	2112286	10.35%	2.016%
50	11.092	3022	3033	3055	rVB	935692	1704307	8.35%	1.627%
51	11.295	3086	3096	3116	rVB	444888	810944	3.97%	0.774%
52	11.471	3141	3151	3165	rVB	2691545	4088308	20.04%	3.903%
53	11.645	3199	3205	3217	rVV	174651	288150	1.41%	0.275%
54	11.748	3220	3237	3245	rVV	316181	571952	2.80%	0.546%
55	11.816	3245	3258	3271	rVB2	773857	1435200	7.03%	1.370%
56	11.899	3271	3284	3298	rBV	1267638	2007838	9.84%	1.917%
57	12.098	3335	3346	3351	rBV2	476959	771447	3.78%	0.736%
58	12.127	3351	3355	3364	rVB	472833	653714	3.20%	0.624%
59	12.192	3364	3375	3396	rVB5	597063	1249123	6.12%	1.192%
60	12.378	3425	3433	3449	rVB	282723	453790	2.22%	0.433%
61	12.468	3450	3461	3465	rBV	316049	460847	2.26%	0.440%
62	12.500	3466	3471	3481	rVV	392763	579827	2.84%	0.553%
63	12.574	3485	3494	3501	rVV	425053	640407	3.14%	0.611%
64	12.616	3501	3507	3517	rVB	141859	214320	1.05%	0.205%
65	12.687	3518	3529	3549	rBV4	479456	1109615	5.44%	1.059%
66	12.880	3579	3589	3598	rVB3	183343	324704	1.59%	0.310%
67	12.976	3599	3619	3627	rBV2	197122	407411	2.00%	0.389%
68	13.031	3627	3636	3652	rVB	343765	589685	2.89%	0.563%
69	13.114	3652	3662	3675	rBV6	113651	282736	1.39%	0.270%
70	13.295	3712	3718	3726	rVB3	156355	234939	1.15%	0.224%
71	13.455	3758	3768	3782	rVB3	714521	1332490	6.53%	1.272%
72	13.629	3814	3822	3846	rVB4	131784	260162	1.27%	0.248%
73	13.841	3879	3888	3900	rBV5	139398	265089	1.30%	0.253%
74	13.944	3915	3920	3935	rVB2	190414	331283	1.62%	0.316%
75	14.089	3950	3965	3987	rVB	485280	896112	4.39%	0.855%
76	15.227	4305	4319	4330	rBV	203599	327206	1.60%	0.312%

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Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
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_U\Method\82U092821W.M
Title : SW846 8260

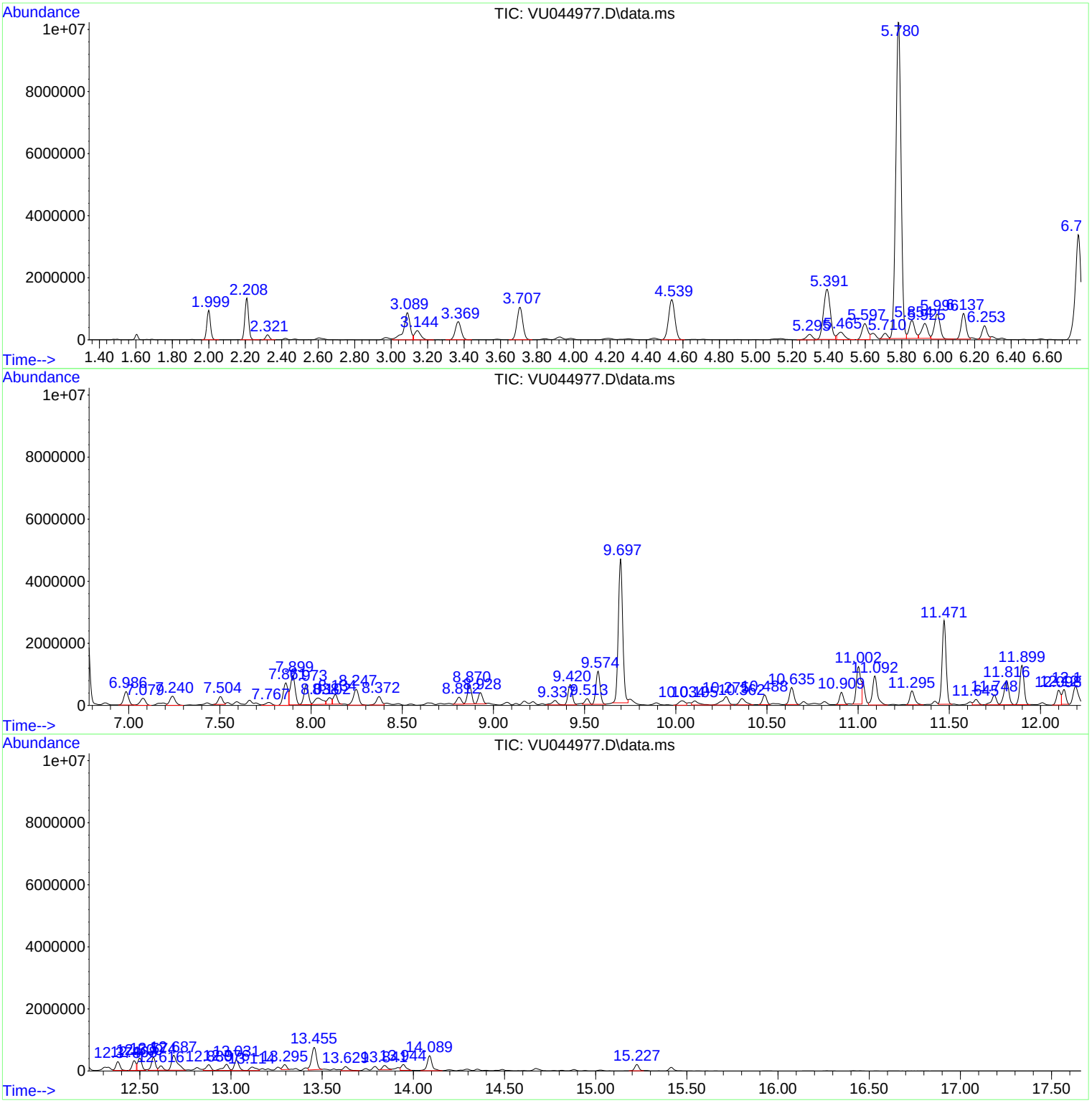
Sum of corrected areas: 104761166

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
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ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

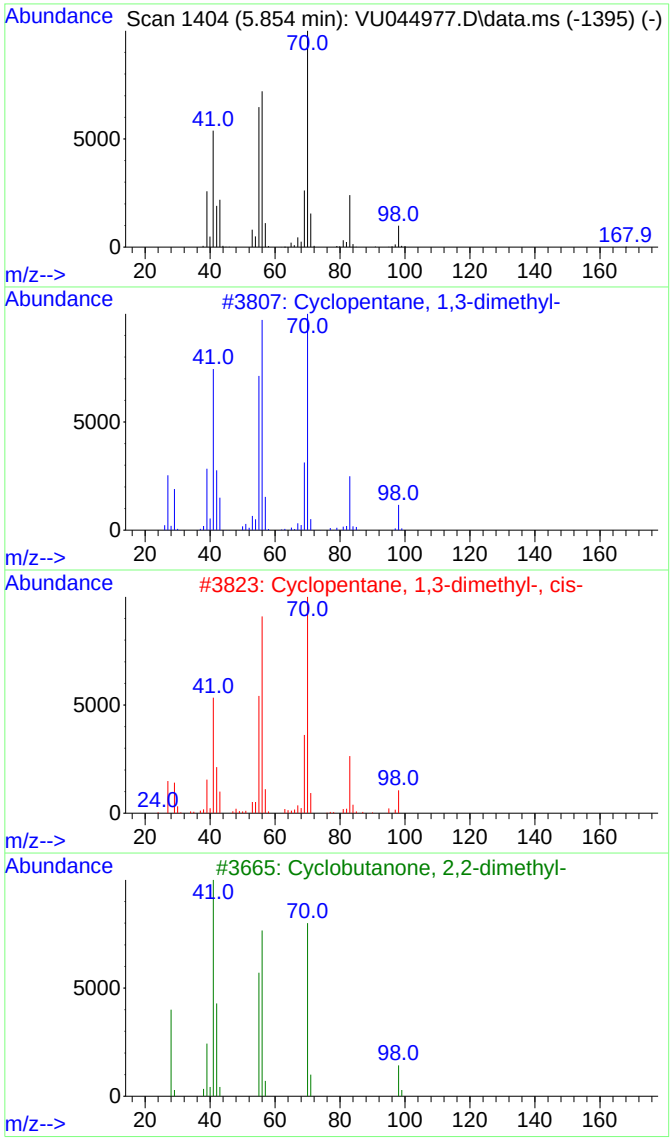
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

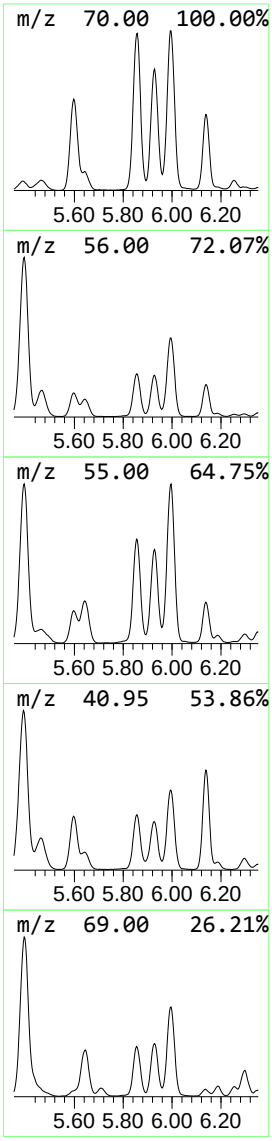
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Cyclopentane, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.854	77.65 ug/l	1252610	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	95
2	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	90
3	Cyclobutanone, 2,2-dimethyl-	98	C6H10O	001192-14-9	72
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	64
5	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	58





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

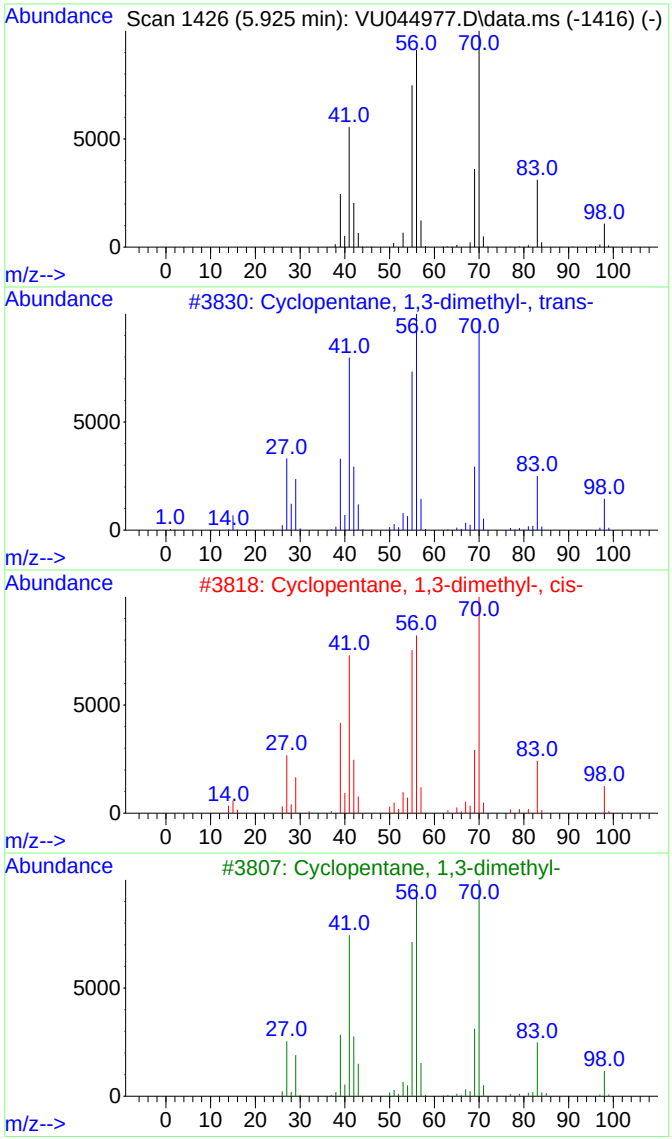
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

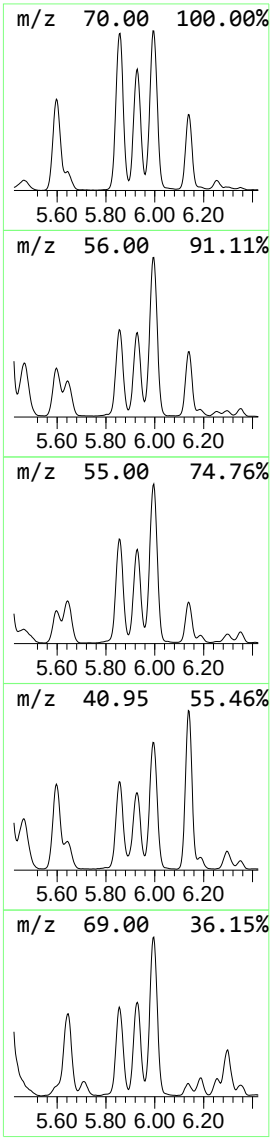
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Cyclopentane, 1,3-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.925	66.22 ug/l	1068120	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, trans-	98	C7H14	001759-58-6	94
2	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	91
3	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	91
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	90
5	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

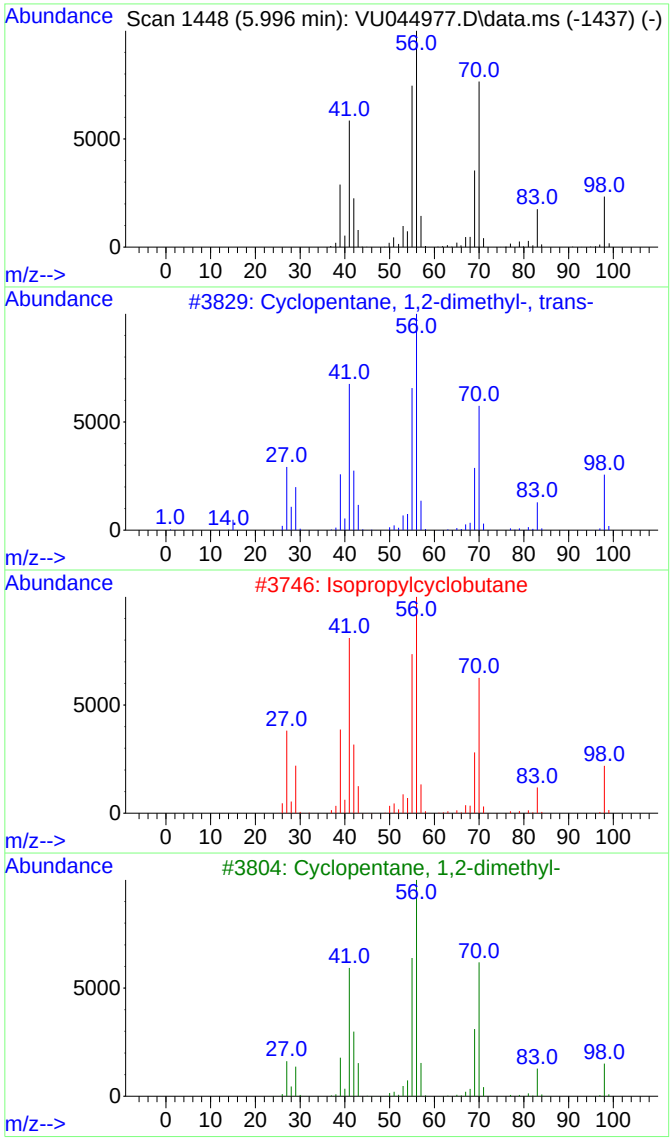
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

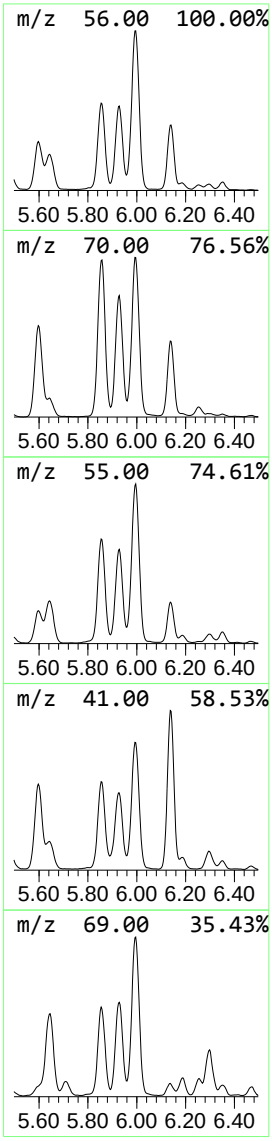
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Cyclopentane, 1,2-dimethyl-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.996	108.61 ug/l	1752000	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	94
2	Isopropylcyclobutane	98	C7H14	000872-56-0	94
3	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	91
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	87
5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	86





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Instrument :
MSVOA_U
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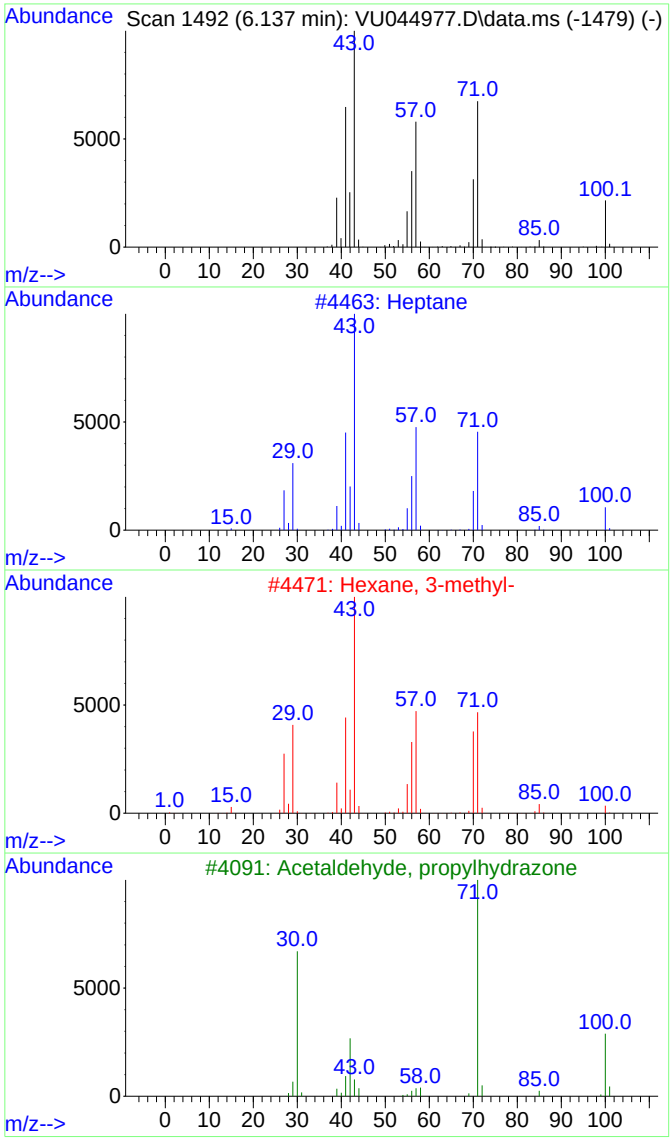
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

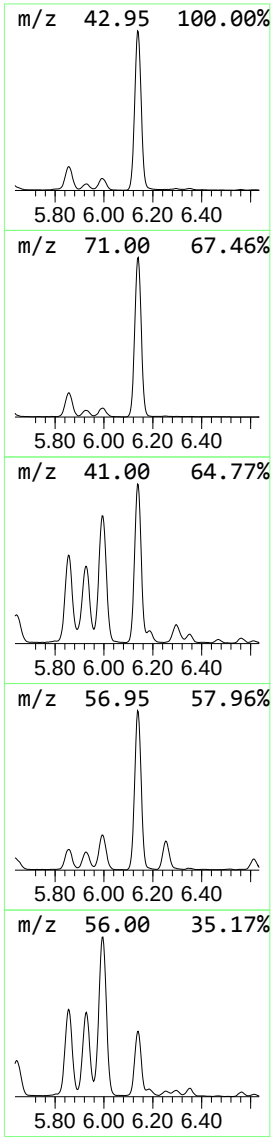
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Heptane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.137	95.61 ug/l	1542240	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane	100	C7H16	000142-82-5	91
2	Hexane, 3-methyl-	100	C7H16	000589-34-4	59
3	Acetaldehyde, propylhydrazone	100	C5H12N2	007422-88-0	47
4	3-Hexanone	100	C6H12O	000589-38-8	46
5	Pentane, 2-methyl-	86	C6H14	000107-83-5	37





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Operator : SY/MD
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Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

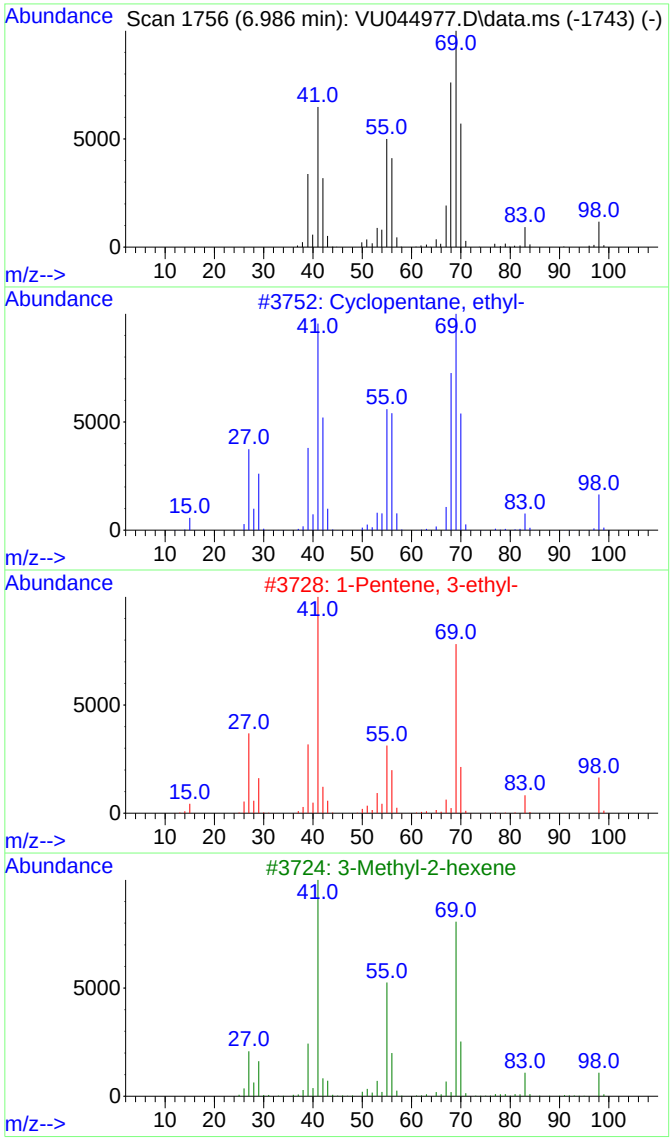
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

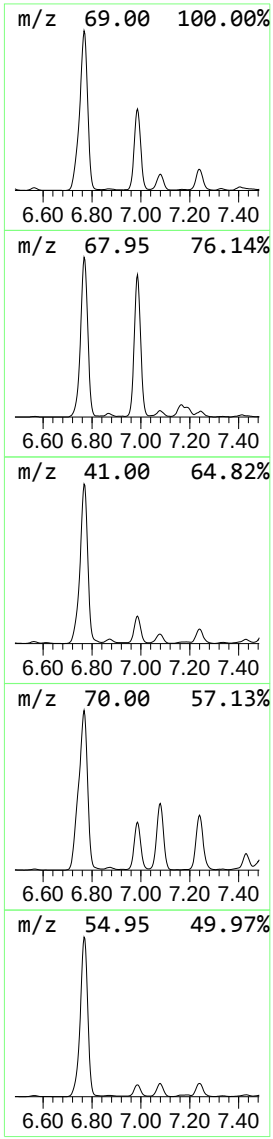
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, ethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.986	48.38 ug/l	780397	1,4-Difluorobenzene	6.253

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, ethyl-	98	C7H14	001640-89-7	81
2	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	43
3	3-Methyl-2-hexene	98	C7H14	017618-77-8	38
4	Cyclopropane, 1,1-diethyl-	98	C7H14	001003-19-6	38
5	Butanedinitrile, 2,3-diamino-2,3...	138	C6H10N4	082929-43-9	35





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Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

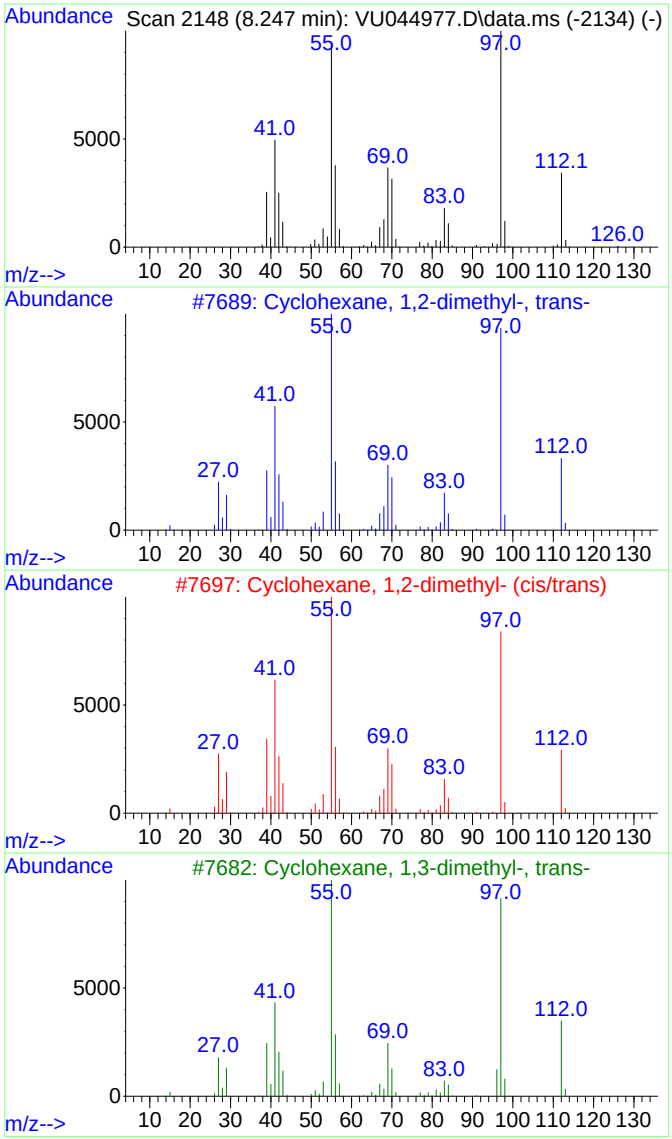
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

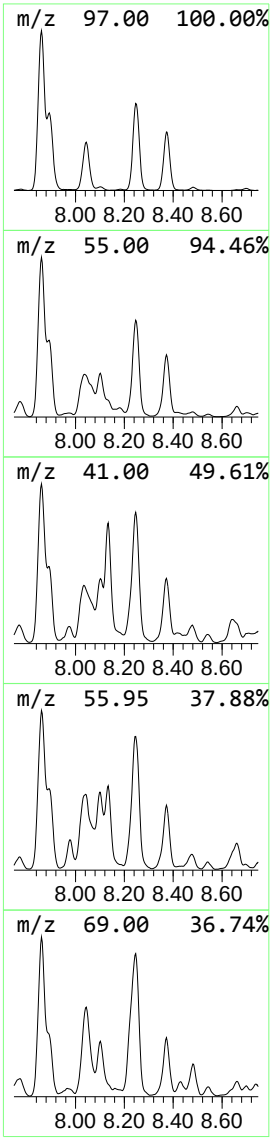
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.247	48.83 ug/l	1037150	Chlorobenzene-d5	9.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	97
2	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	95
3	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
4	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	83
5	Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

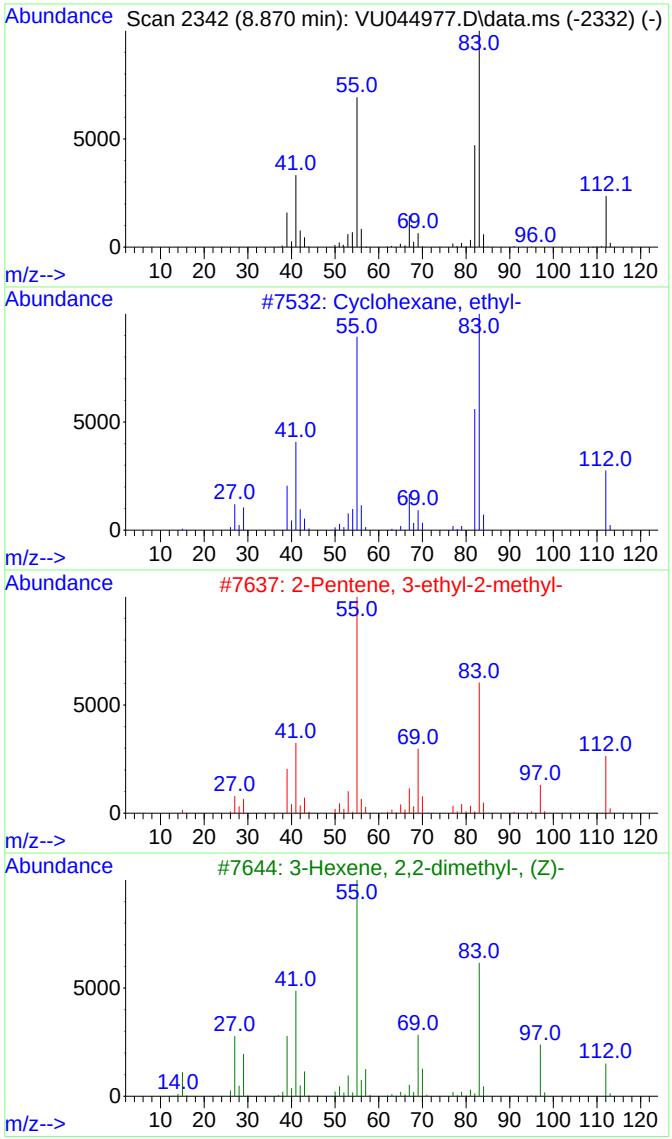
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

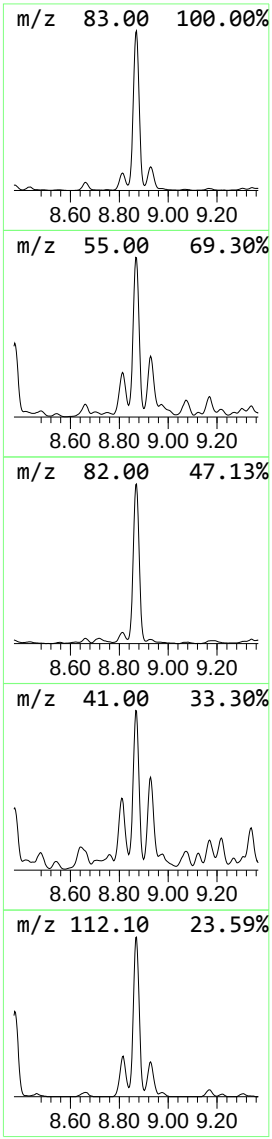
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclohexane, ethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.870	44.59 ug/l	947030	Chlorobenzene-d5	9.420

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	97
2	2-Pentene, 3-ethyl-2-methyl-	112	C8H16	019780-67-7	58
3	3-Hexene, 2,2-dimethyl-, (Z)-	112	C8H16	000690-92-6	53
4	Oxalic acid, cyclohexyl dodecyl ...	340	C20H36O4	1000309-31-4	53
5	2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

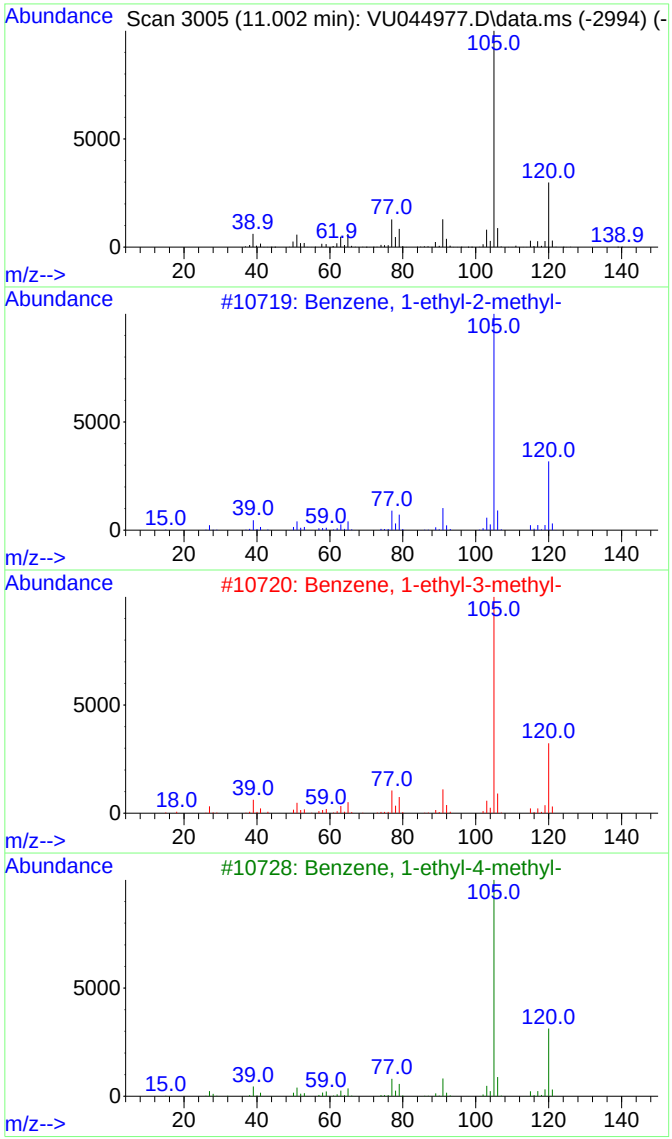
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

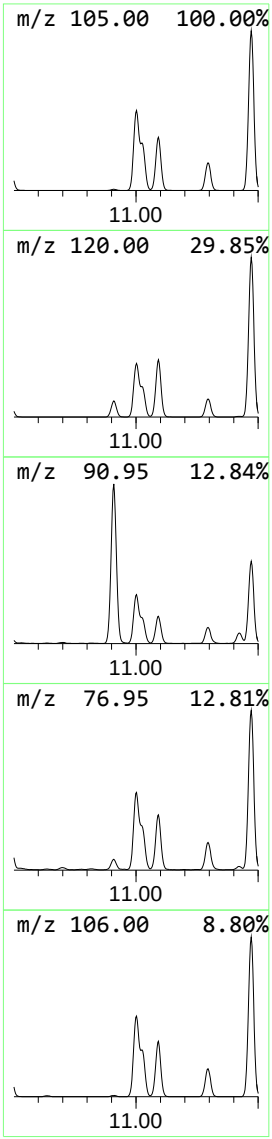
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.002	73.59 ug/l	2112290	1,4-Dichlorobenzene-d4	11.816

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-3-methyl-	120	C9H12	000620-14-4	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

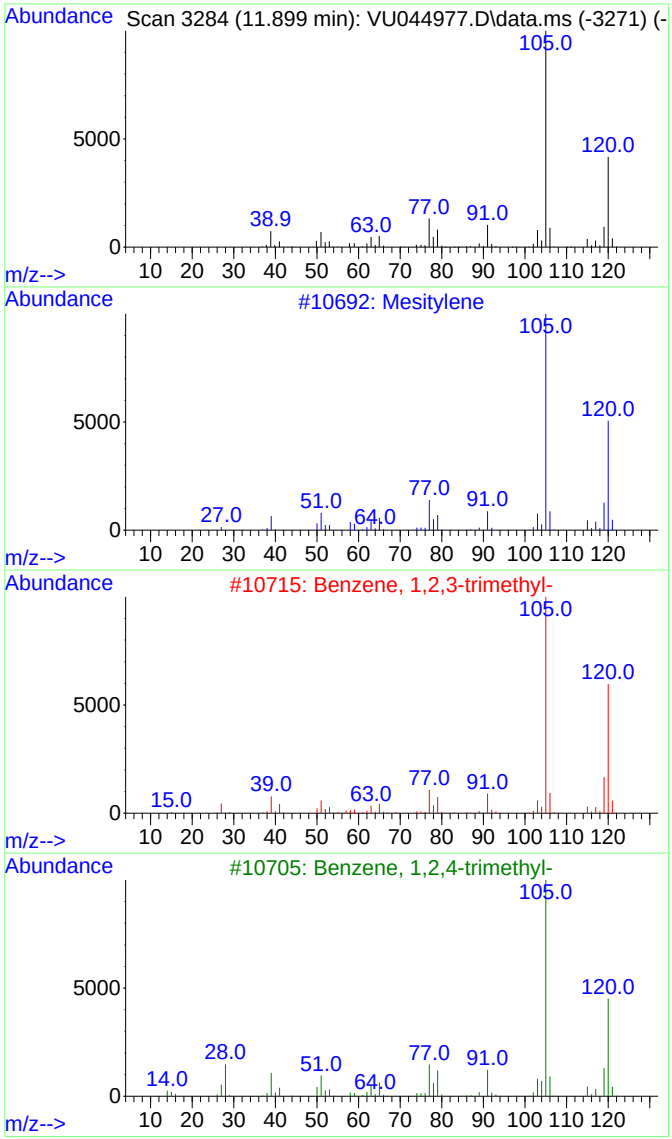
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

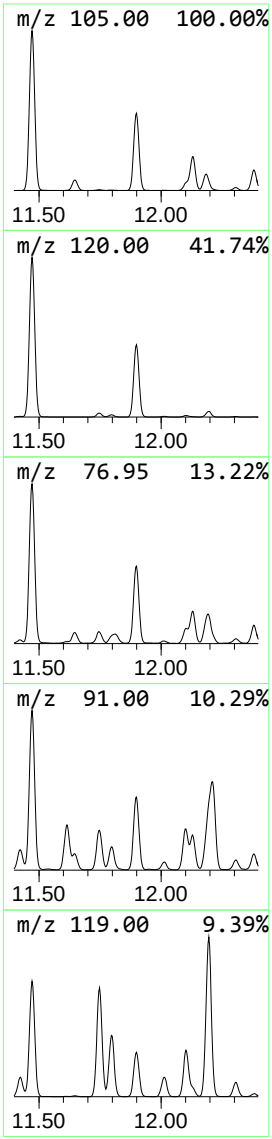
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	69.95 ug/l	2007840	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Mesitylene	120	C9H12	000108-67-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	90
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

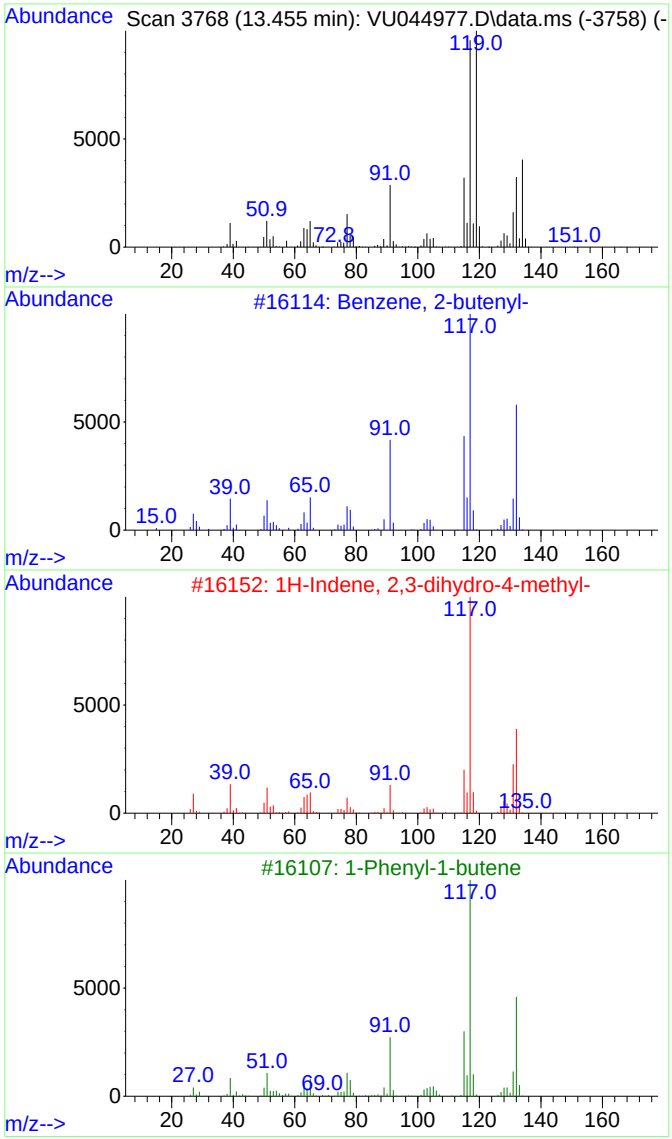
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

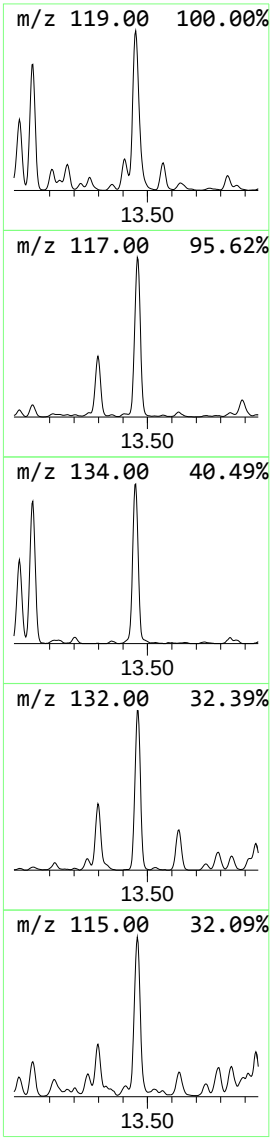
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-butenyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.455	46.42 ug/l	1332490	1,4-Dichlorobenzene-d4	11.816

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	86
3	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	64





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092821\
Data File : VU044977.D
Acq On : 29 Sep 2021 06:04
Operator : SY/MD
Sample : M3952-05
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 52 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MW-3

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.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
Cyclopentane, 1...	5.854	77.7	ug/l	1252610	2	6.253	806530	50.0
Cyclopentane, 1...	5.925	66.2	ug/l	1068120	2	6.253	806530	50.0
Cyclopentane, 1...	5.996	108.6	ug/l	1752000	2	6.253	806530	50.0
Heptane	6.137	95.6	ug/l	1542240	2	6.253	806530	50.0
Cyclopentane, e...	6.986	48.4	ug/l	780397	2	6.253	806530	50.0
Cyclohexane, 1,...	8.247	48.8	ug/l	1037150	3	9.420	1061910	50.0
Cyclohexane, et...	8.870	44.6	ug/l	947030	3	9.420	1061910	50.0
Benzene, 1-ethy...	11.002	73.6	ug/l	2112290	4	11.816	1435200	50.0
Benzene, 1,2,3-...	11.899	70.0	ug/l	2007840	4	11.816	1435200	50.0
Benzene, 2-bute...	13.455	46.4	ug/l	1332490	4	11.816	1435200	50.0