LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SUP-UST-03-(8-10)ME

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % 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: VU044996.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.292	1212	1229	1241	rBV	184370	401439	13.80%	0.631%
2	5.372	1241	1254	1271	rVV	393287	858188	29.50%	1.349%
3	5.584	1306	1320	1331	rBV2	37935	84246	2.90%	0.132%
4	5.703	1343	1357	1374	rVB	232492	472204	16.23%	0.742%
5	5.986	1433	1445	1462	rVB3	32490	70133	2.41%	0.110%
6	6.131	1476	1490	1504	rBV	51555	97080	3.34%	0.153%
7	6.250	1511	1527	1560	rBV	506420	1000556	34.40%	1.572%
8	6.761	1664	1686	1698	rBV	205632	454666	15.63%	0.715%
9	7.237	1822	1834	1852	rVB2	34110	77409	2.66%	0.122%
10	7.497	1902	1915	1923	rBV	118205	207839	7.15%	0.327%
11	7.658	1956	1965	1974	rBV2	64353	102476	3.52%	0.161%
12	7.858	2011	2027	2031	rBV2	131012	229790	7.90%	0.361%
13	7.899	2031	2040	2056	rVB	829315	1500342	51.58%	2.358%
14	8.034	2067	2082	2095	rBV5	40532	133789	4.60%	0.210%
15	8.243	2133	2147	2161	rVB3	109425	212959	7.32%	0.335%
16	8.369	2174	2186	2195	rBV2	50797	95090	3.27%	0.149%
17	8.636	2250	2269	2282	rBV2	126971	239187	8.22%	0.376%
18	8.761	2297	2308	2314	rBV	56158	100589	3.46%	0.158%
19	8.806	2314	2322	2331	rVB5	69325	119915	4.12%	0.188%
20	8.867	2332	2341	2349	rBV	150455	226434	7.78%	0.356%
21	8.925	2349	2359	2370	rBV	368142	622890	21.41%	0.979%
22	9.073	2393	2405	2413	rBV3	61312	115360	3.97%	0.181%
23	9.121	2413	2420	2427	rVV	101987	157253	5.41%	0.247%
24	9.169	2427	2435	2442	rVV4	120363	221211	7.60%	0.348%
25	9.214	2442	2449	2459	rVB2	168795	270786	9.31%	0.426%
26	9.337	2477	2487	2502	rVB	267327	430960	14.82%	0.677%
27	9.423	2502	2514	2526	rBV	758300	1296853	44.58%	2.038%
28	9.565	2549	2558	2572	rVB2	70167	129360	4.45%	0.203%
29	9.671	2573	2591	2597	rBV6	154785	391403	13.46%	0.615%
30	9.716	2598	2605	2613	rVV5	188038	329294	11.32%	0.518%
31	9.764	2613	2620	2645	rVB2	129101	302927	10.41%	0.476%
32	9.883	2645	2657	2670	rBV6	69679	170533	5.86%	0.268%
33	10.025	2689	2701	2714	rBV6	218179	590688	20.31%	0.928%
34	10.124	2723	2732	2742	rBV4	114870	233519	8.03%	0.367%
35	10.253	2755	2772	2786	rBV4	556805	1224777	42.11%	1.925%
36	10.359	2795	2805	2813	rBV4	377298	717899	24.68%	1.128%

LSC Area Percent Report

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Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SUP-UST-03-(8-10)ME

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 1 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

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Peak separation: 5

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

37	10.485	2832	2844	2852	rVB4	77351	175774	6.04%	0.276%
38	10.558	2852	2867	2875	rBV7	162720	419786	14.43%	0.660%
39	10.636	2876	2891	2904	rVB3	853784	1759848	60.50%	2.766%
40	10.700	2904	2911	2918	rBV2	78459	124654	4.29%	0.196%
41	10.761	2923	2930	2936	rVV2	129989	198613	6.83%	0.312%
42	10.812	2937	2946	2963	rVB4	548752	989869	34.03%	1.556%
43	10.906	2965	2975	2982	rBV2	129195	238162	8.19%	0.374%
44	10.951	2983	2989	3002	rVV3	132722	249147	8.57%	0.392%
45	11.021	3003	3011	3017	rVV5	151976	286347	9.84%	0.450%
46	11.066	3018	3025	3035	rVV2	358096	651641	22.40%	1.024%
47	11.131	3037	3045	3055	rVB5	252338	527538	18.14%	0.829%
48	11.189	3056	3063	3074	rBV5	108722	184724	6.35%	0.290%
49	11.250	3075	3082	3087	rBV5	70475	107622	3.70%	0.169%
50	11.314	3097	3102	3114	rVB7	227681	388032	13.34%	0.610%
51	11.382	3114	3123	3127	rBV4	177590	290913	10.00%	0.457%
52	11.420	3128	3135	3143	rVB	594412	799525	27.49%	1.257%
53	11.465	3143	3149	3155	rBV6	132345	215517	7.41%	0.339%
54	11.578	3176	3184	3189	rBV2	127255	214459	7.37%	0.337%
55	11.645	3199	3205	3216	rVB4	406865	677015	23.27%	1.064%
56	11.716	3218	3227	3234	rBV2	419546	667922	22.96%	1.050%
57	11.819	3249	3259	3271	rBV2	958808	1757353	60.41%	2.762%
58	11.877	3272	3277	3284	rVB3	148582	182688	6.28%	0.287%
59	11.992	3306	3313	3326	rVB6	351091	645645	22.20%	1.015%
60	12.102	3337	3347	3358	rBV2	499339	930203	31.98%	1.462%
61	12.189	3366	3374	3378	rBV4	417893	643732	22.13%	1.012%
62	12.388	3430	3436	3450	rVB6	328122	701713	24.12%	1.103%
63	12.574	3488	3494	3501	rVB	399850	525865	18.08%	0.826%
64	12.684	3517	3528	3549	rVB6	447993	1267770	43.58%	1.992%
65	12.841	3571	3577	3593	rVB6	355422	806297	27.72%	1.267%
66	12.935	3597	3606	3612	rBV5	430519	856875	29.46%	1.347%
67	13.057	3638	3644	3652	rVB4	337082	504179	17.33%	0.792%
68	13.111	3654	3661	3676	rVB4	259649	539426	18.54%	0.848%
69	13.259	3702	3707	3714	rBV4	176604	233180	8.02%	0.366%
70	13.455	3759	3768	3785	rVB3	1484053	2908852	100.00%	4.571%
71	13.565	3797	3802	3806	rBV	164460	194630	6.69%	0.306%
72	13.738	3848	3856	3863	rBV3	338321	568195	19.53%	0.893%
73	13.790	3865	3872	3879	rVV	353801	538293	18.51%	0.846%
74	13.841	3880	3888	3897	rVV5	590397	986861	33.93%	1.551%
75	13.947	3916	3921	3933	rVB2	384843	679520	23.36%	1.068%
76	14.124	3969	3976	3985	rVB5	378843	565473	19.44%	0.889%

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ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SUP-UST-03-(8-10)ME

Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
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Title : SW846 8260

77	14.353	4042	4047	4053	rBV4	165096	227729	7.83%	0.358%
78	14.488	4082	4089	4096	rBV2	421246	632776	21.75%	0.994%
79	14.671	4137	4146	4159	rBV4	841997	1652939	56.82%	2.598%
80	14.816	4184	4191	4204	rBV3	396189	716557	24.63%	1.126%
81	14.883	4205	4212	4221	rVB2	710879	1112071	38.23%	1.748%
82	14.944	4225	4231	4246	rVB5	326107	654734	22.51%	1.029%
83	15.025	4247	4256	4278	rBV10	482246	1326721	45.61%	2.085%
84	15.224	4311	4318	4329	rVB	1289745	2081263	71.55%	3.271%
85	15.349	4351	4357	4363	rBV6	209005	336105	11.55%	0.528%
86	15.414	4369	4377	4389	rVB2	1209697	2103076	72.30%	3.305%
87	16.153	4599	4607	4613	rBV2	612683	937100	32.22%	1.473%
88	16.266	4633	4642	4652	rBV	1632852	2631526	90.47%	4.136%
89	16.414	4679	4688	4694	rBV	1738583	2696700	92.71%	4.238%
90	16.452	4694	4700	4716	rVB	1292122	2140994	73.60%	3.365%
91	16.533	4718	4725	4738	rVB3	263396	491646	16.90%	0.773%
92	16.642	4745	4759	4780	rVB	558202	1649423	56.70%	2.592%
93	16.790	4797	4805	4826	rVB	361433	681915	23.44%	1.072%
94	16.883	4828	4834	4845	rVB3	164940	252651	8.69%	0.397%
95	17.121	4901	4908	4922	rVB	287984	471229	16.20%	0.741%
96	17.227	4935	4941	4951	rVB5	120011	196784	6.77%	0.309%
97	17.330	4967	4973	4981	rVB2	297004	448052	15.40%	0.704%
98	17.378	4982	4988	5018	rVB	258581	550395	18.92%	0.865%
99	17.542	5032	5039	5050	rVB	204918	342013	11.76%	0.537%
100	17.610	5052	5060	5071	rVV	118298	202225	6.95%	0.318%

Sum of corrected areas: 63630526

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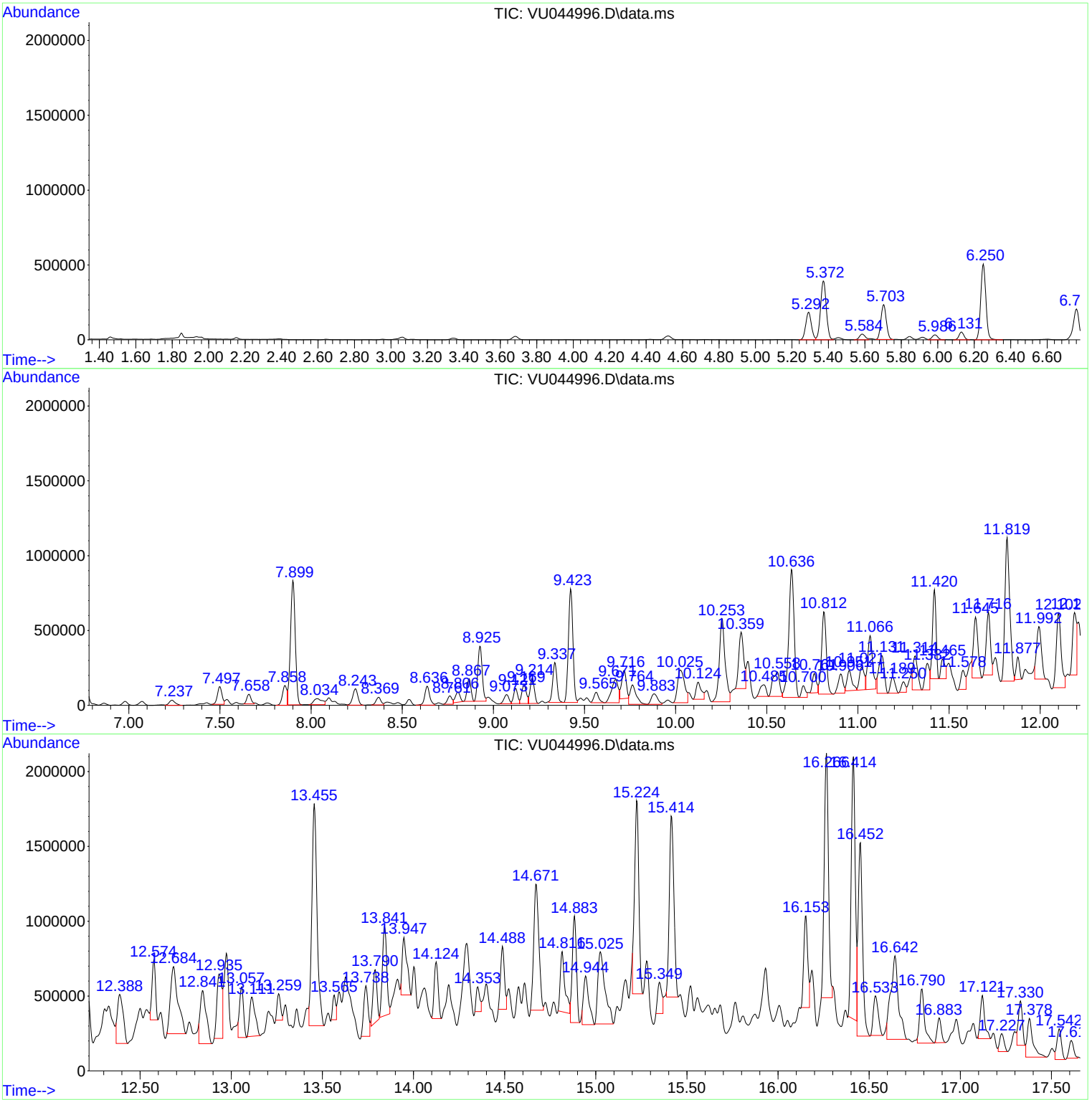
Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

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\VU092921\
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Acq On : 29 Sep 2021 16:28
Operator : SY/MD
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Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

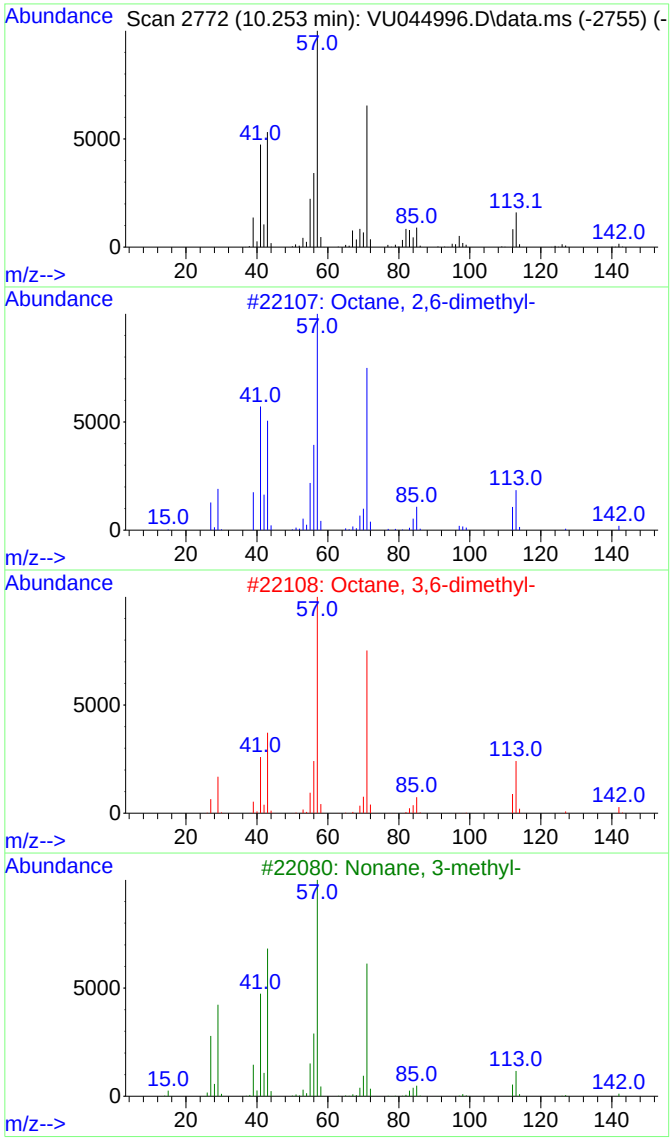
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

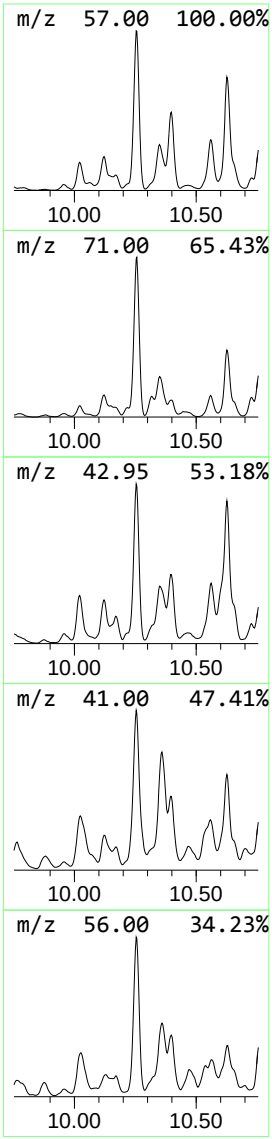
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.253	47.22 ug/l	1224780	Chlorobenzene-d5	9.423

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	94
2	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	91
3	Nonane, 3-methyl-	142	C10H22	005911-04-6	90
4	Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	53
5	Butane, 2,2-dimethyl-	86	C6H14	000075-83-2	53





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Operator : SY/MD
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ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

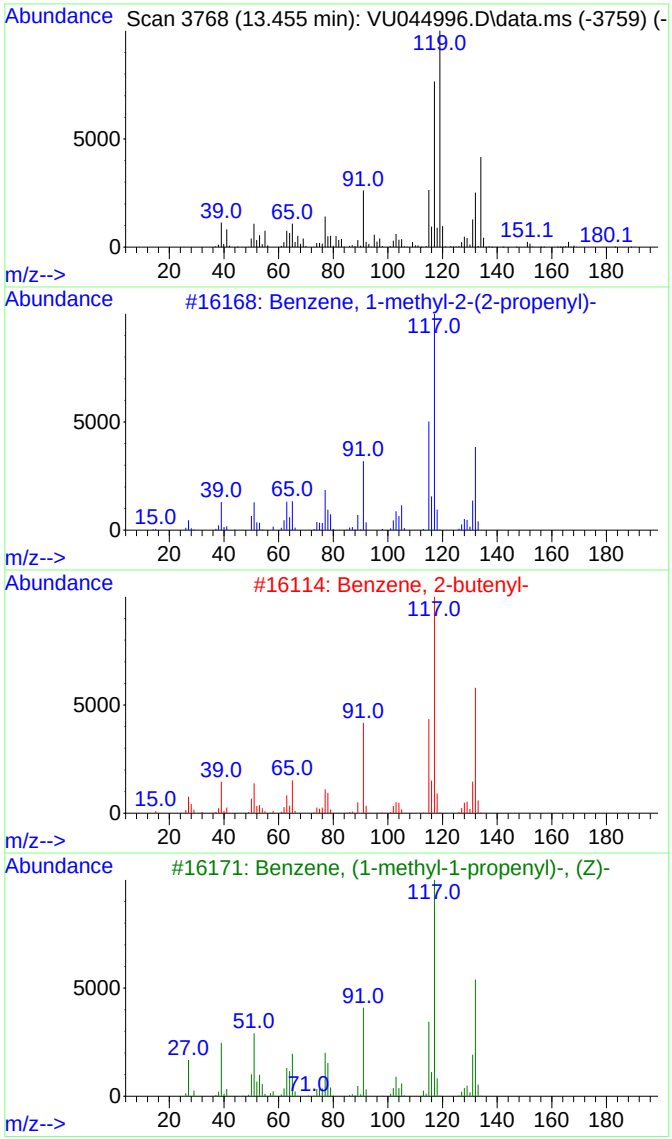
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

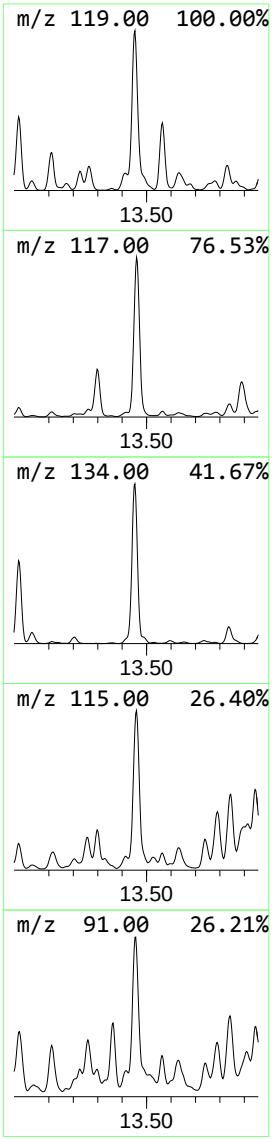
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.456	82.76 ug/l	2908850	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
3	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	64
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	64
5	p-Cymene	134	C10H14	000099-87-6	55





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Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

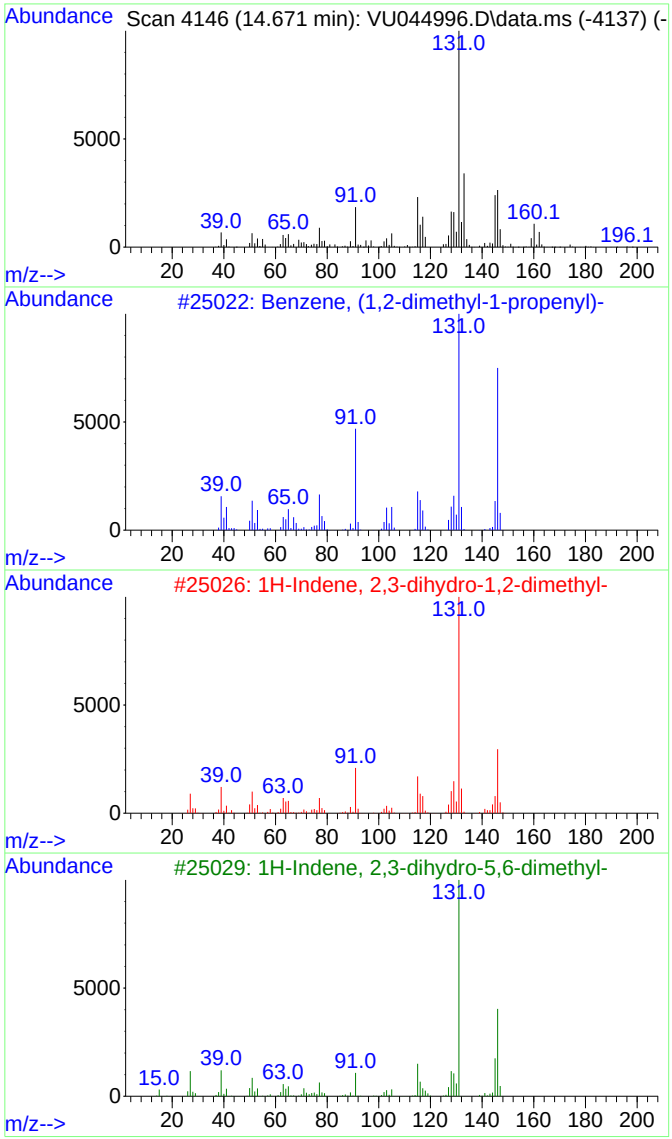
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

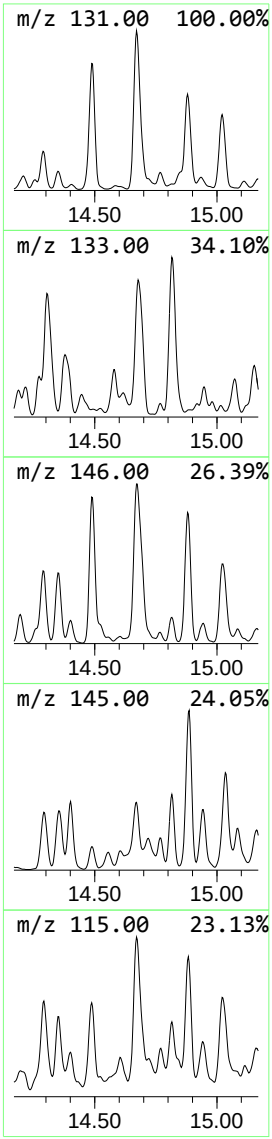
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, (1,2-dimethyl-1-pr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.671	47.03 ug/l	1652940	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	93
2	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	76
3	1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	70
4	Naphthalene, 6-ethyl-1,2,3,4-tet...	160	C12H16	022531-20-0	70
5	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	64





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Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

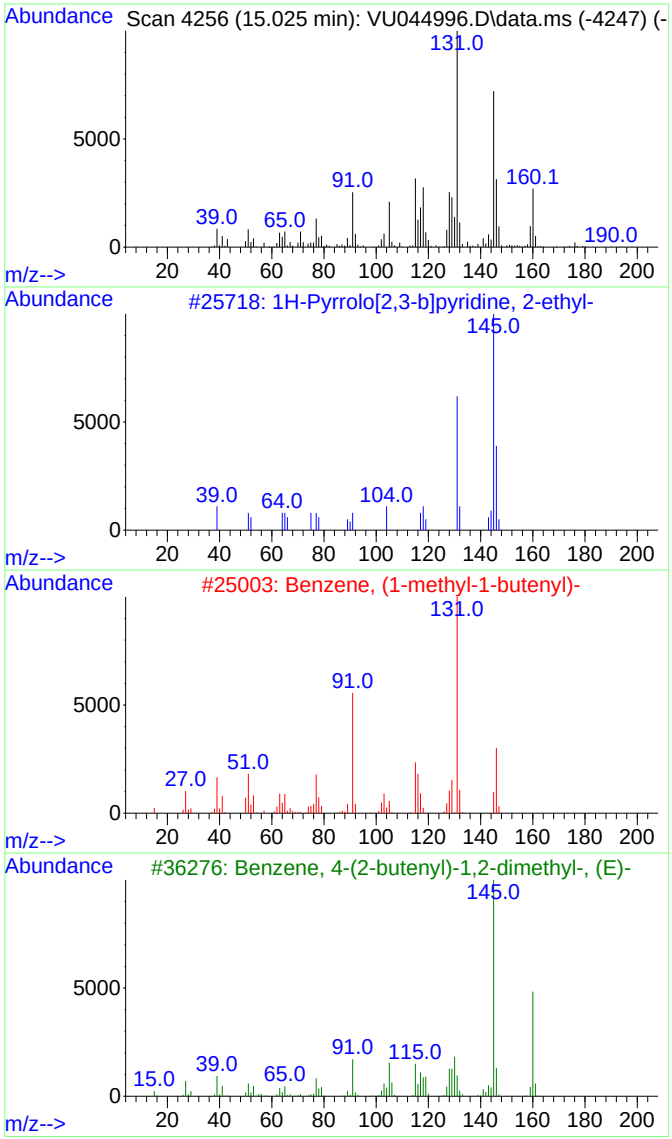
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

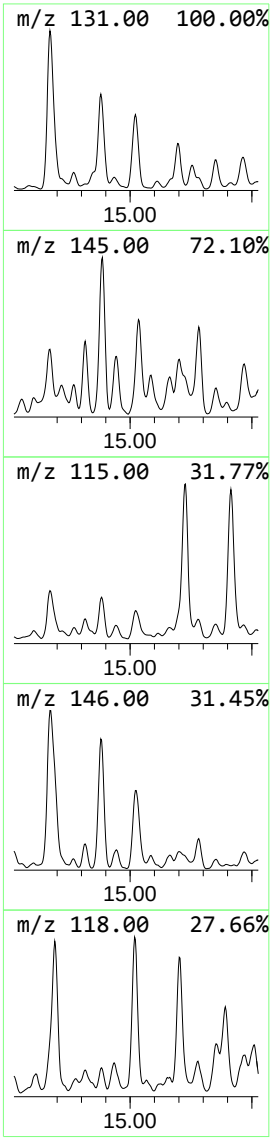
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 1H-Pyrrolo[2,3-b]pyridine, ... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.025	37.75 ug/l	1326720	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Pyrrolo[2,3-b]pyridine, 2-ethyl-	146	C9H10N2	023612-49-9	52
2	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	49
3	Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	49
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	46
5	Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	46





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ClientSampleId :
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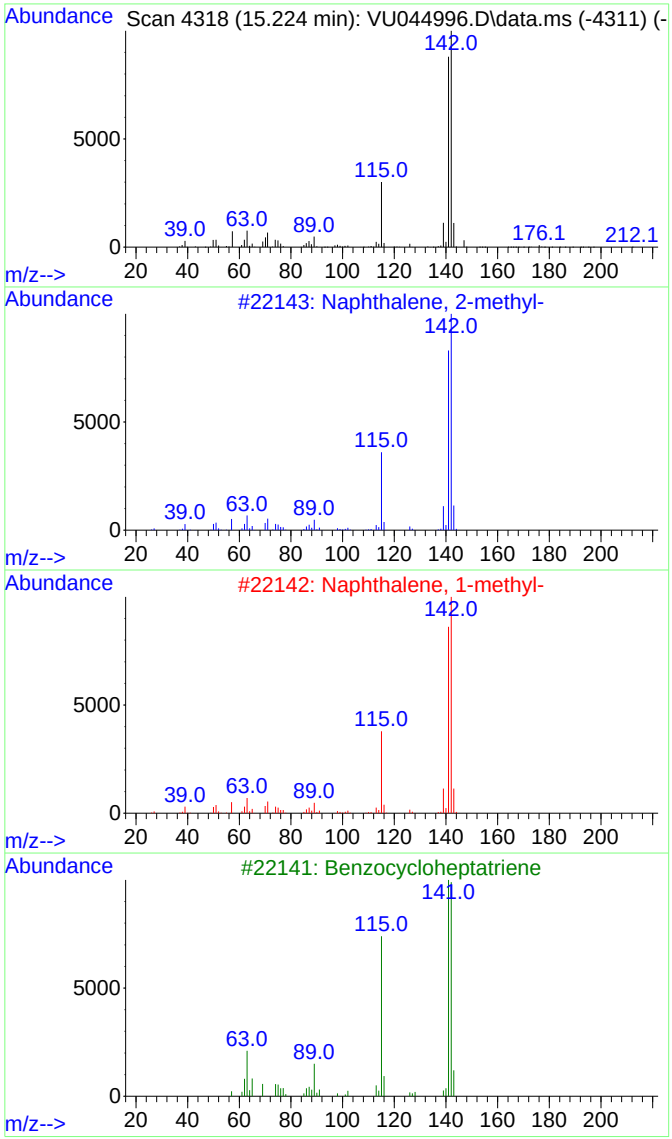
Quant Title : SW846 8260

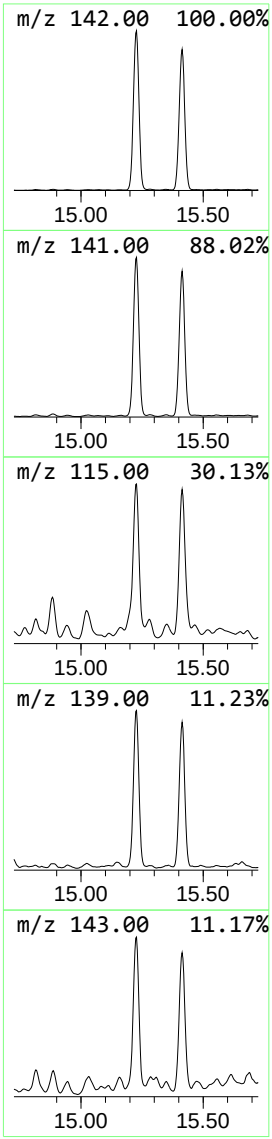
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Naphthalene, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.224	59.22 ug/l	2081260	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	94
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	94
3	Benzocycloheptatriene	142	C11H10	000264-09-5	91
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	91
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

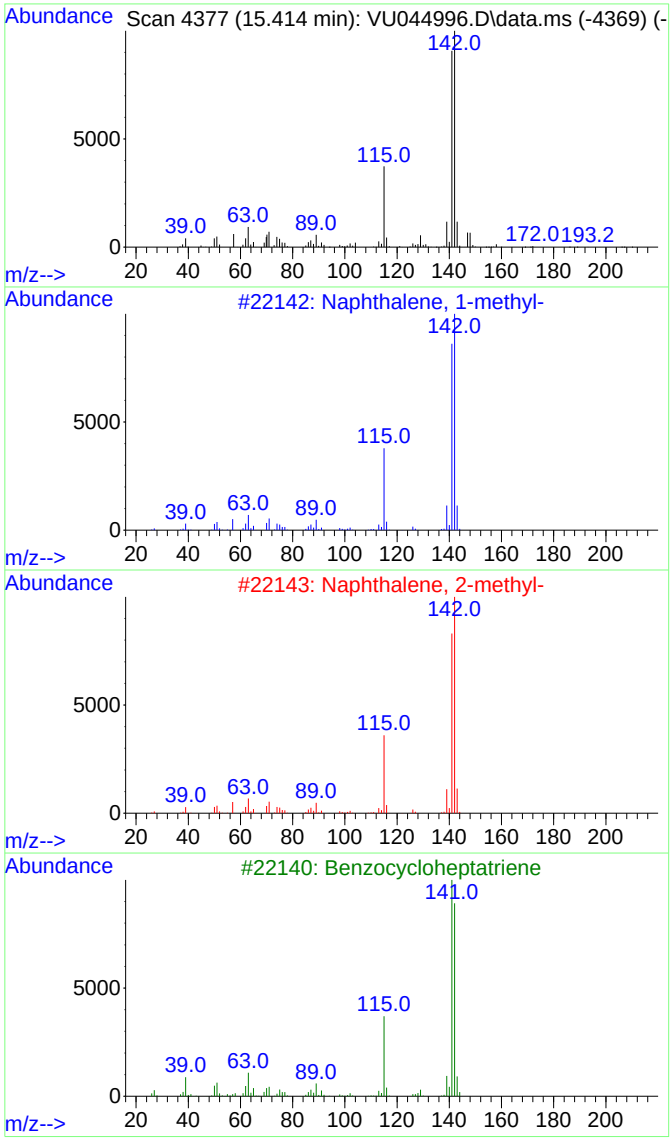
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

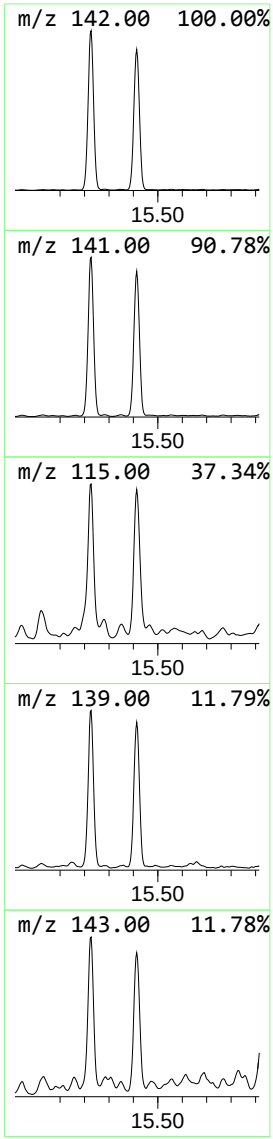
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Naphthalene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.414	59.84 ug/l	2103080	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	94
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	93
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

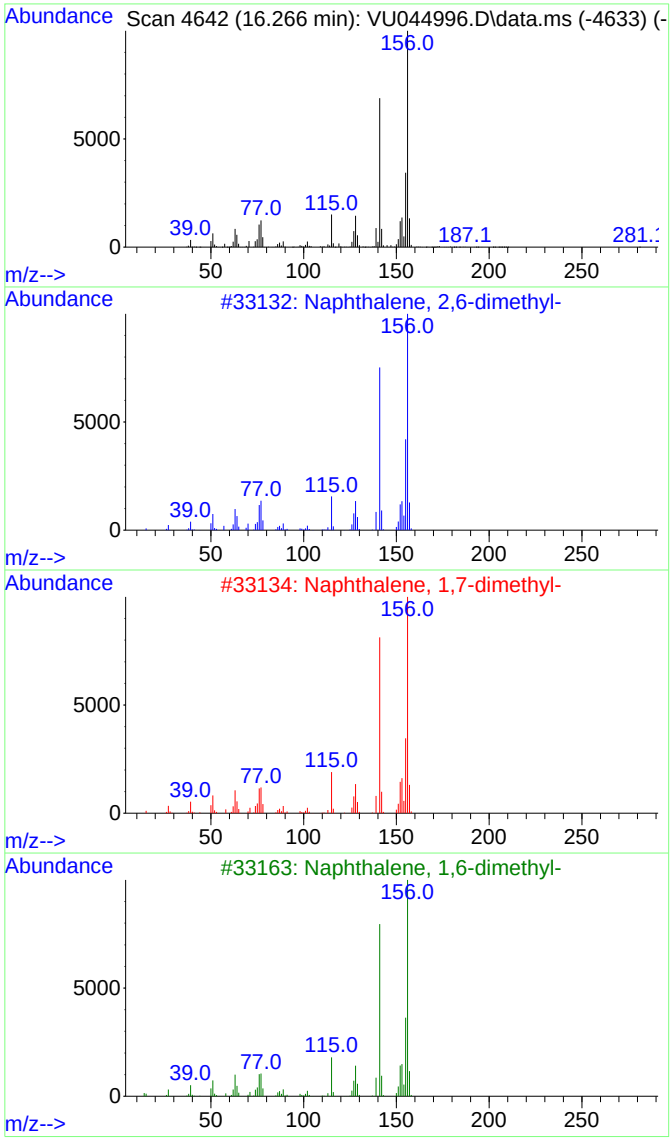
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

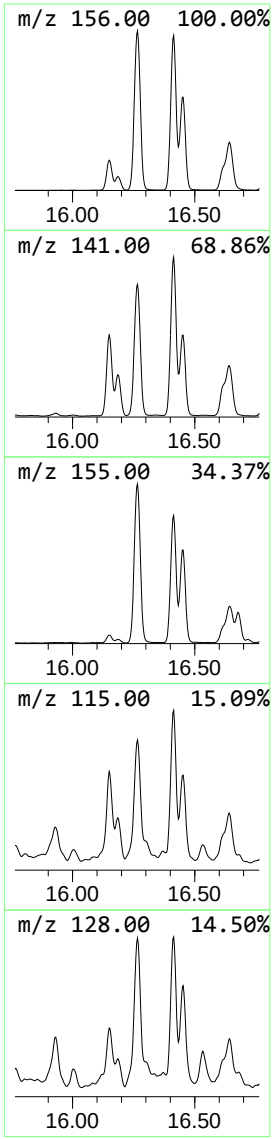
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Naphthalene, 2,6-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.266	74.87 ug/l	2631530	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	98
2	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
3	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
4	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

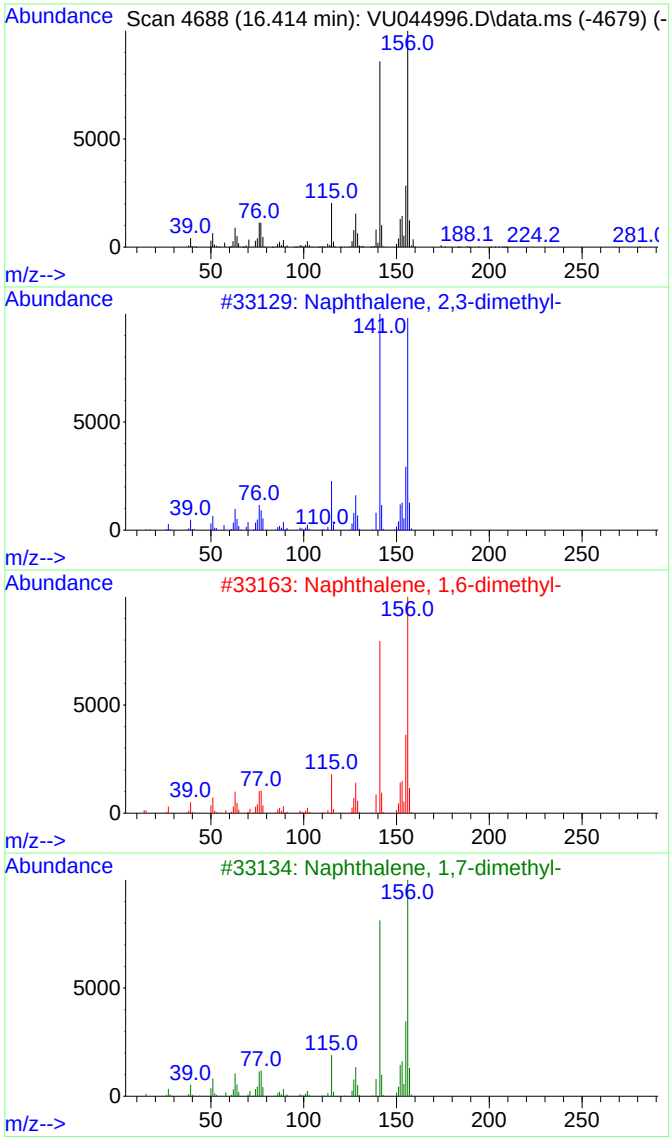
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

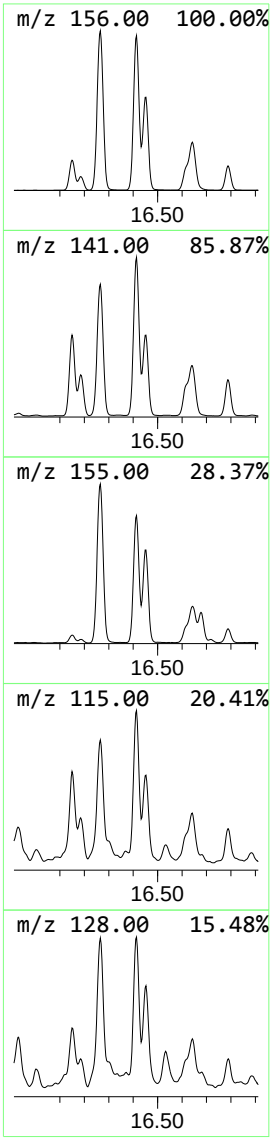
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.413	76.73 ug/l	2696700	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
2	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	98
4	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
5	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

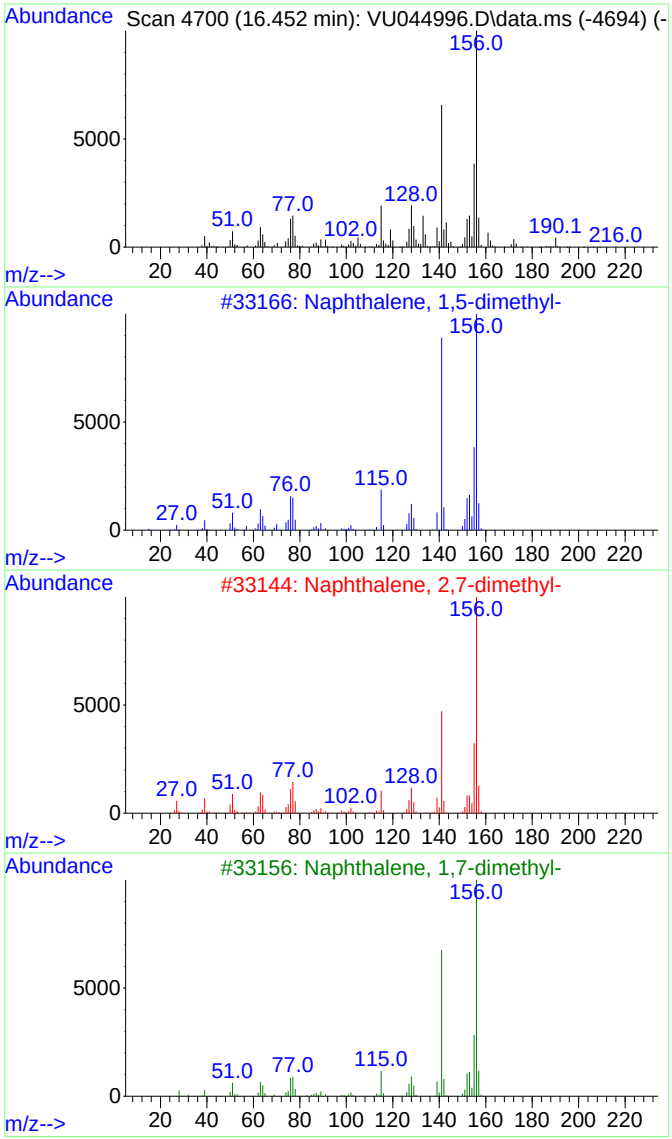
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

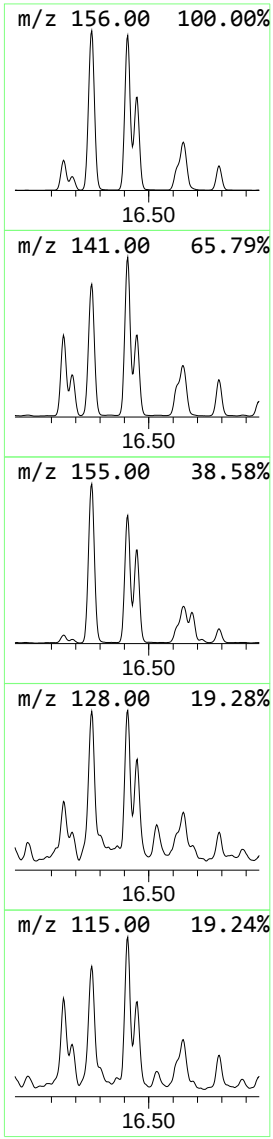
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Naphthalene, 1,5-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.452	60.92 ug/l	2140990	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
2	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
4	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5	Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
SUP-UST-03-(8-10)ME

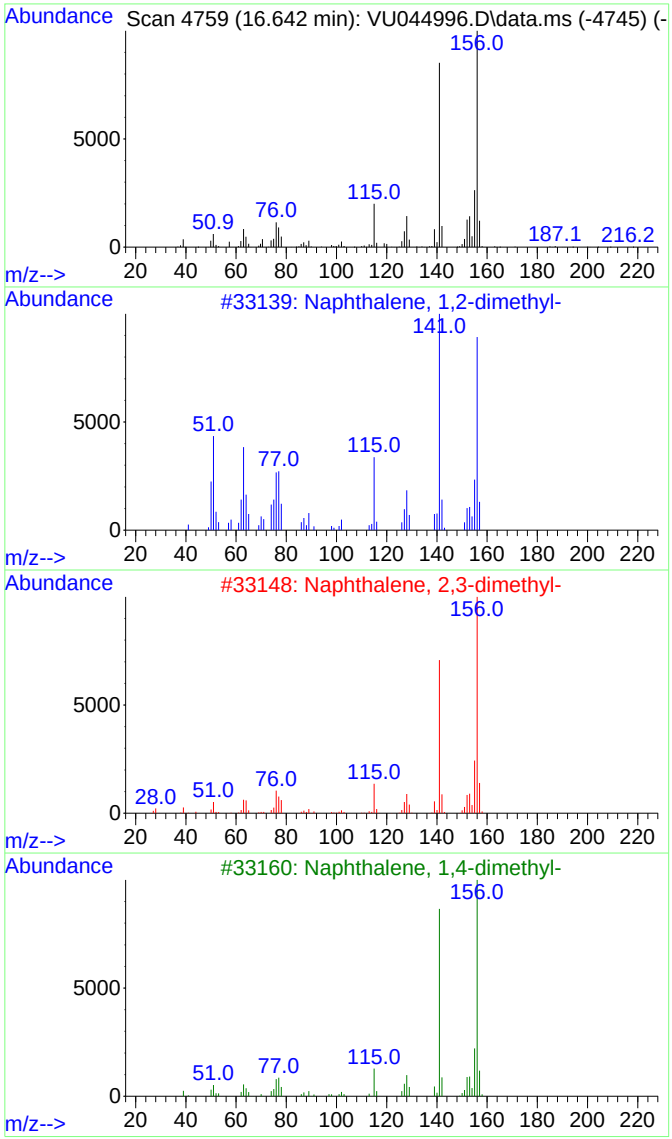
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U092821W.M
Quant Title : SW846 8260

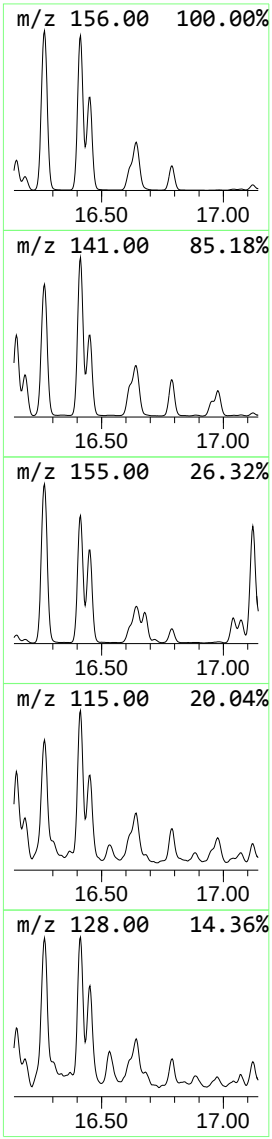
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Naphthalene, 1,2-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.642	46.93 ug/l	1649420	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
3	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4	Naphthalene, 1,8-dimethyl-	156	C12H12	000569-41-5	95
5	Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092921\
Data File : VU044996.D
Acq On : 29 Sep 2021 16:28
Operator : SY/MD
Sample : M3889-02ME
Misc : 7.04g/5mL/100uL/5.00mL/MSVOA_U/MEOH
ALS Vial : 13 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : SUP-UST-03-(8-10)ME

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
Octane, 2,6-dim...	10.253	47.2	ug/l	1224780	3	9.423	1296850	50.0
Benzene, 1-meth...	13.456	82.8	ug/l	2908850	4	11.819	1757350	50.0
Benzene, (1,2-d...	14.671	47.0	ug/l	1652940	4	11.819	1757350	50.0
1H-Pyrrolo[2,3-...	15.025	37.7	ug/l	1326720	4	11.819	1757350	50.0
Naphthalene, 2-...	15.224	59.2	ug/l	2081260	4	11.819	1757350	50.0
Naphthalene, 1-...	15.414	59.8	ug/l	2103080	4	11.819	1757350	50.0
Naphthalene, 2,...	16.266	74.9	ug/l	2631530	4	11.819	1757350	50.0
Naphthalene, 2,...	16.413	76.7	ug/l	2696700	4	11.819	1757350	50.0
Naphthalene, 1,...	16.452	60.9	ug/l	2140990	4	11.819	1757350	50.0
Naphthalene, 1,...	16.642	46.9	ug/l	1649420	4	11.819	1757350	50.0