LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMULM092922WMA.M
Title : VOC Analysis

Signal : TIC: VU051019.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.597	74	80	97	rVB	221233	263500	14.95%	1.768%
2	1.906	170	176	198	rVB	214084	285123	16.17%	1.913%
3	2.565	369	381	396	rBV	341863	553135	31.38%	3.711%
4	3.044	518	530	538	rBV5	2215	4086	0.23%	0.027%
5	3.353	616	626	638	rBV3	6387	12194	0.69%	0.082%
6	3.433	648	651	655	rBV2	167	152	0.01%	0.001%
7	3.462	655	660	665	rVB2	347	356	0.02%	0.002%
8	3.533	678	682	686	rBV2	255	240	0.01%	0.002%
9	3.565	689	692	696	rBV2	277	250	0.01%	0.002%
10	3.642	713	716	717	rBV2	283	143	0.01%	0.001%
11	3.742	744	747	750	rBV2	201	153	0.01%	0.001%
12	4.060	842	846	848	rBV	322	191	0.01%	0.001%
13	4.096	853	857	861	rBV	353	342	0.02%	0.002%
14	4.150	869	874	881	rBV2	656	859	0.05%	0.006%
15	4.182	881	884	888	rVB	425	370	0.02%	0.002%
16	4.314	920	925	927	rBV	409	327	0.02%	0.002%
17	4.369	939	942	945	rBV2	257	145	0.01%	0.001%
18	4.562	999	1002	1005	rBV2	271	176	0.01%	0.001%
19	4.623	1005	1021	1050	rBV	172309	474743	26.93%	3.185%
20	5.060	1142	1157	1181	rBV	320799	702615	39.86%	4.713%
21	5.433	1270	1273	1276	rBV	499	239	0.01%	0.002%
22	5.449	1276	1278	1282	rVB2	251	163	0.01%	0.001%
23	5.539	1302	1306	1307	rBV	220	137	0.01%	0.001%
24	5.594	1320	1323	1326	rBV2	295	197	0.01%	0.001%
25	5.722	1342	1363	1375	rBV2	639345	1762840	100.00%	11.826%
26	6.005	1449	1451	1454	rBV2	573	379	0.02%	0.003%
27	6.089	1473	1477	1480	rBV4	474	452	0.03%	0.003%
28	6.247	1510	1526	1550	rBV	487479	930579	52.79%	6.243%
29	6.533	1603	1615	1633	rVB8	17320	40236	2.28%	0.270%
30	6.603	1633	1637	1647	rBV4	655	1004	0.06%	0.007%
31	6.690	1649	1664	1682	rBV	415437	814161	46.18%	5.462%
32	6.925	1735	1737	1741	rVB	264	158	0.01%	0.001%
33	6.944	1741	1743	1745	rBV2	258	143	0.01%	0.001%
34	7.047	1771	1775	1777	rBV	302	226	0.01%	0.002%
35	7.105	1789	1793	1798	rVB2	307	260	0.01%	0.002%
36	7.137	1798	1803	1804	rBV	246	167	0.01%	0.001%

LSC Area Percent Report

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Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMULM092922WMA.M
Title : VOC Analysis

37	7.185	1811	1818	1821	rBV4	1397	1525	0.09%	0.010%
38	7.304	1852	1855	1859	rBV	283	212	0.01%	0.001%
39	7.565	1923	1936	1956	rBV	183319	330736	18.76%	2.219%
40	7.735	1987	1989	1991	rVB	508	222	0.01%	0.001%
41	7.806	2005	2011	2012	rBV2	360	298	0.02%	0.002%
42	7.896	2025	2039	2054	rBV	700718	1198301	67.98%	8.039%
43	8.179	2115	2127	2149	rBV	131016	229473	13.02%	1.539%
44	8.327	2171	2173	2175	rBV3	470	261	0.01%	0.002%
45	8.472	2211	2218	2219	rBV4	673	681	0.04%	0.005%
46	8.552	2237	2243	2252	rVB5	3310	4955	0.28%	0.033%
47	8.636	2256	2269	2309	rBV3	431642	946809	53.71%	6.352%
48	8.822	2324	2327	2332	rBV2	755	450	0.03%	0.003%
49	9.005	2381	2384	2387	rBV2	304	231	0.01%	0.002%
50	9.079	2404	2407	2408	rBV	215	144	0.01%	0.001%
51	9.144	2422	2427	2430	rVB3	365	346	0.02%	0.002%
52	9.172	2430	2436	2439	rBV2	346	356	0.02%	0.002%
53	9.327	2482	2484	2487	rVB2	643	298	0.02%	0.002%
54	9.417	2498	2512	2534	rBV	691324	1154567	65.49%	7.745%
55	9.568	2548	2559	2580	rVB	210830	341296	19.36%	2.290%
56	9.693	2588	2598	2617	rVB	55364	93168	5.29%	0.625%
57	9.870	2644	2653	2662	rVB6	2040	4088	0.23%	0.027%
58	9.983	2685	2688	2691	rBV	351	260	0.01%	0.002%
59	10.053	2704	2710	2713	rVV4	1103	1336	0.08%	0.009%
60	10.098	2714	2724	2743	rVB	66942	112172	6.36%	0.752%
61	10.234	2764	2766	2771	rBV2	321	240	0.01%	0.002%
62	10.333	2792	2797	2801	rBV3	944	943	0.05%	0.006%
63	10.433	2824	2828	2831	rBV	387	241	0.01%	0.002%
64	10.481	2832	2843	2853	rBV	25150	38824	2.20%	0.260%
65	10.687	2905	2907	2909	rBV	333	168	0.01%	0.001%
66	10.754	2915	2928	2953	rBV	590270	950946	53.94%	6.379%
67	10.902	2968	2974	2981	rVB5	2760	3773	0.21%	0.025%
68	10.996	2996	3003	3008	rBV2	3586	4730	0.27%	0.032%
69	11.086	3026	3031	3037	rBV4	1323	1641	0.09%	0.011%
70	11.195	3058	3065	3066	rBV3	524	512	0.03%	0.003%
71	11.237	3075	3078	3082	rBV2	229	161	0.01%	0.001%
72	11.291	3086	3095	3108	rVV3	14235	25383	1.44%	0.170%
73	11.369	3117	3119	3123	rVB	471	241	0.01%	0.002%
74	11.414	3128	3133	3137	rBV4	723	793	0.04%	0.005%
75	11.520	3164	3166	3170	rVB2	396	204	0.01%	0.001%
76	11.642	3198	3204	3214	rVB7	3756	4720	0.27%	0.032%

LSC Area Percent Report

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Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0 % 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\SFAMULM092922WMA.M
Title : VOC Analysis

77	11.809	3244	3256	3273	rBV	574585	913051	51.79%	6.125%
78	12.005	3311	3317	3325	rVB3	4117	5414	0.31%	0.036%
79	12.092	3335	3344	3360	rVB	65068	106659	6.05%	0.716%
80	12.192	3362	3375	3389	rBV	650218	1054473	59.82%	7.074%
81	12.311	3401	3412	3428	rBV	290335	462160	26.22%	3.100%
82	12.603	3500	3503	3512	rVB6	1460	1885	0.11%	0.013%
83	12.680	3516	3527	3540	rBV4	13824	24230	1.37%	0.163%
84	12.806	3561	3566	3575	rBV5	3410	4554	0.26%	0.031%
85	12.922	3591	3602	3611	rBV	96333	155255	8.81%	1.042%
86	13.021	3626	3633	3641	rBV2	33303	52649	2.99%	0.353%
87	13.323	3720	3727	3739	rVB4	7832	12845	0.73%	0.086%
88	13.520	3778	3788	3799	rBV3	18138	27718	1.57%	0.186%
89	13.626	3811	3821	3836	rVV3	61891	104258	5.91%	0.699%
90	13.696	3840	3843	3844	rBV2	656	384	0.02%	0.003%
91	13.831	3877	3885	3892	rBV4	16266	25254	1.43%	0.169%
92	13.880	3894	3900	3919	rVB2	23937	41179	2.34%	0.276%
93	13.970	3924	3928	3930	rBV4	824	777	0.04%	0.005%
94	14.057	3945	3955	3960	rBV	70620	113195	6.42%	0.759%
95	14.542	4091	4106	4126	rBV	11238	43358	2.46%	0.291%
96	14.680	4137	4149	4162	rBV9	7594	18214	1.03%	0.122%
97	14.802	4177	4187	4196	rBV5	27128	47392	2.69%	0.318%
98	15.690	4460	4463	4470	rBV5	1391	1212	0.07%	0.008%
99	16.182	4610	4616	4624	rVB4	10903	15456	0.88%	0.104%
100	17.095	4890	4900	4925	rBV	202318	368989	20.93%	2.475%

Sum of corrected areas: 14906707

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Data File : VU051019.D
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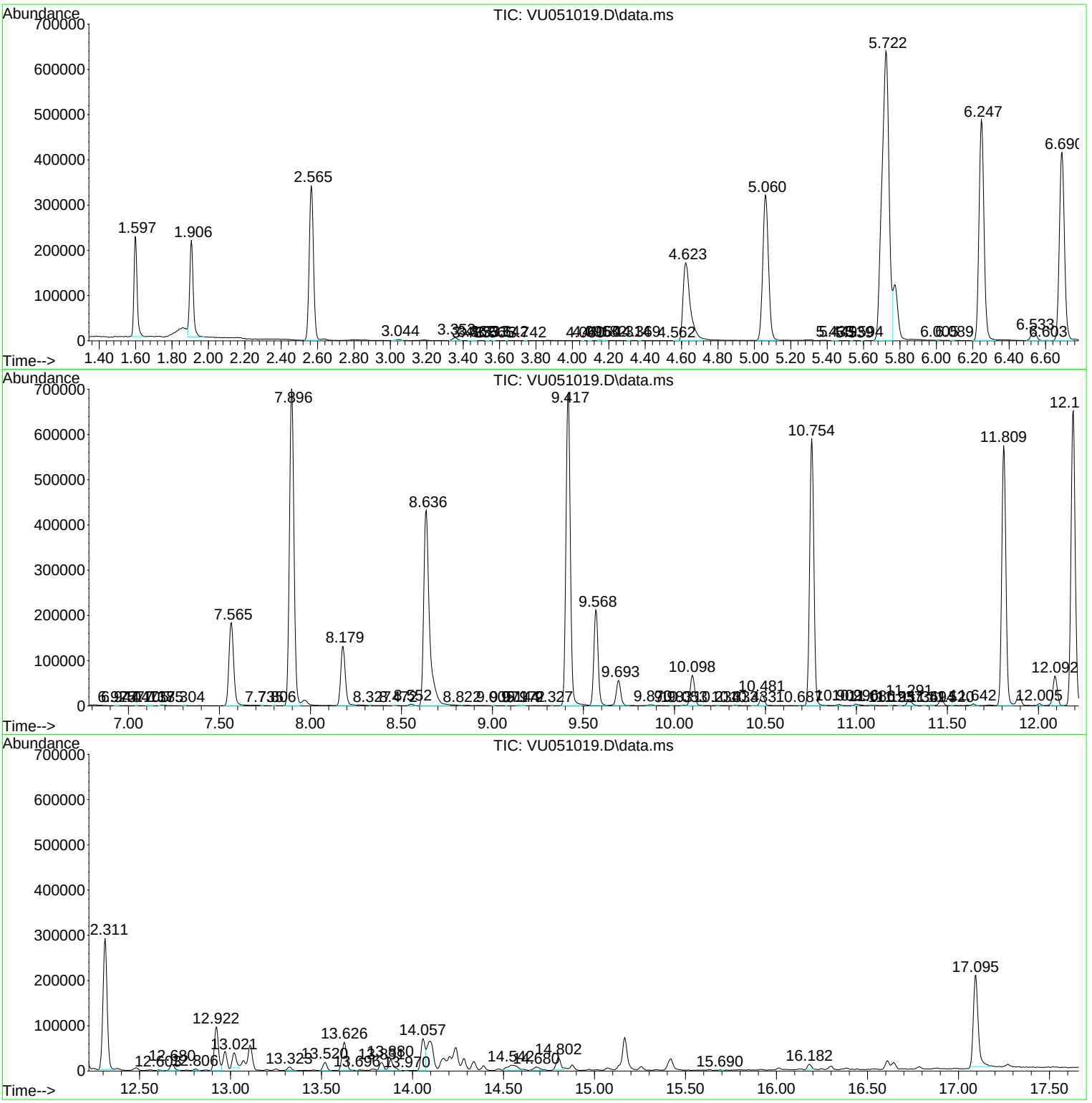
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

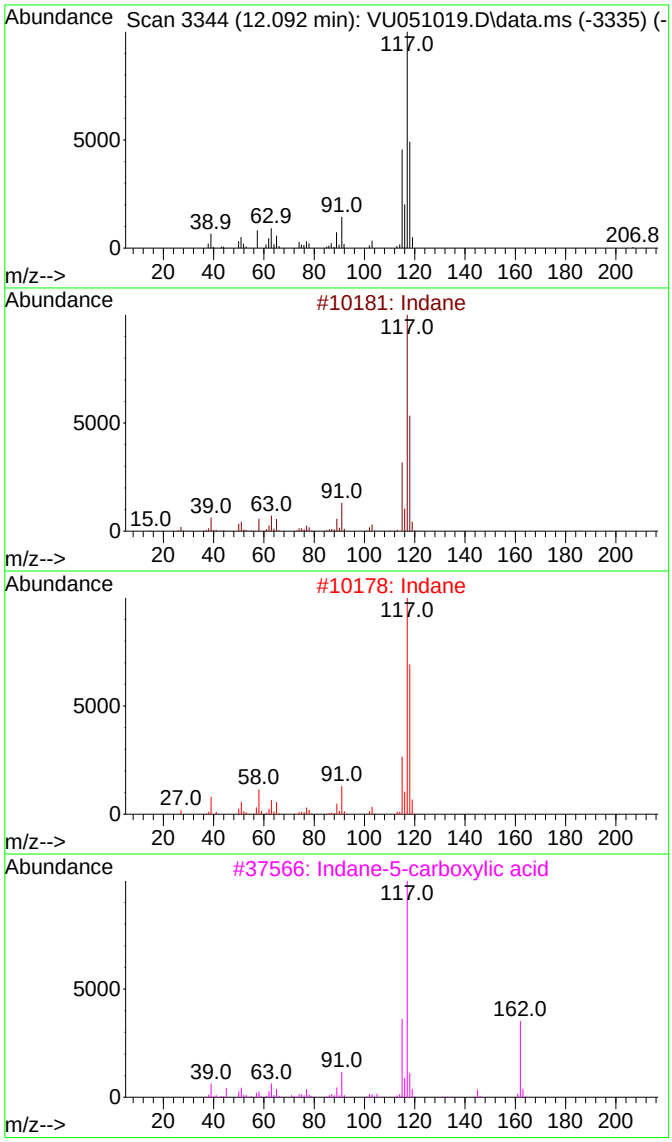
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

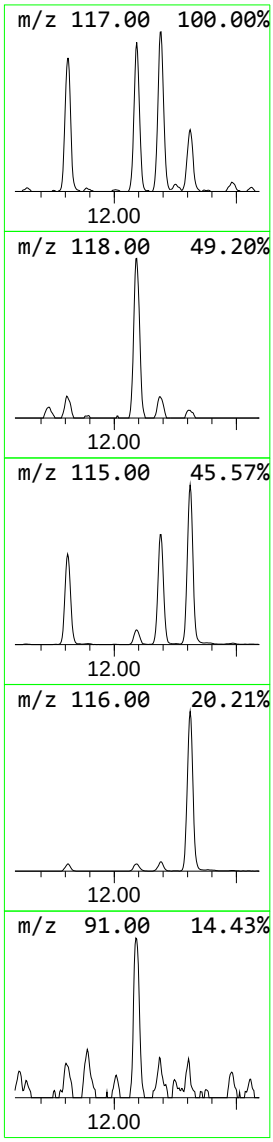
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.092	5.84 ug/L	106659	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Indane	118	C9H10	000496-11-7	64
3	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	59
4	Indane-5-carboxylic acid	162	C10H10O2	065898-38-6	53
5	Benzene, cyclopropyl-	118	C9H10	000873-49-4	52





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

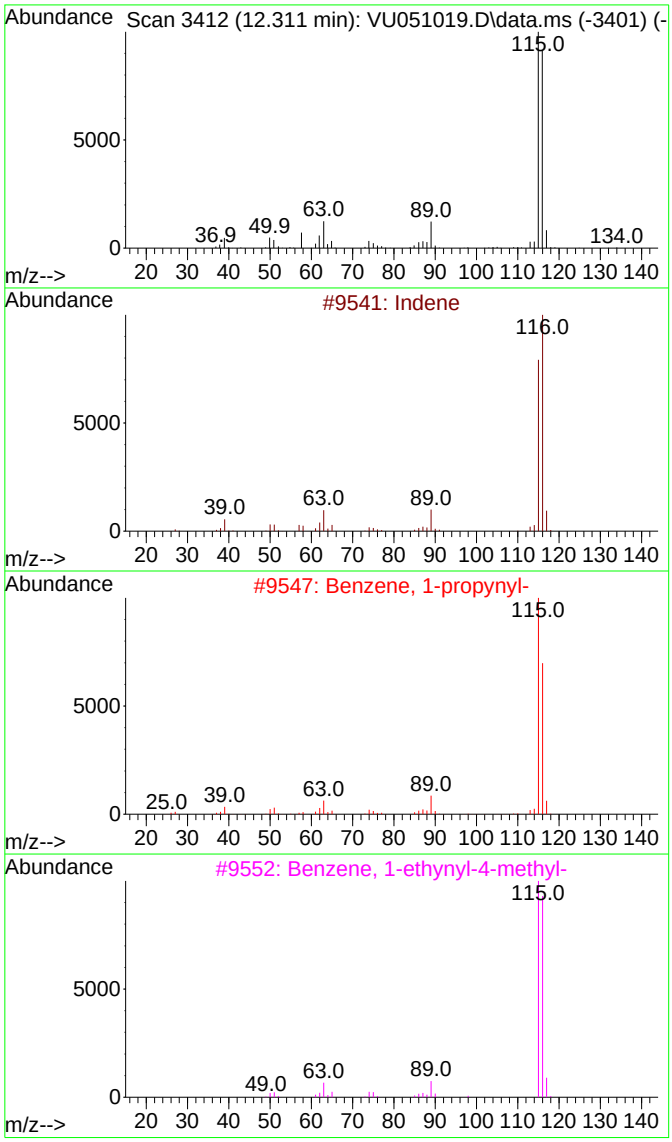
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

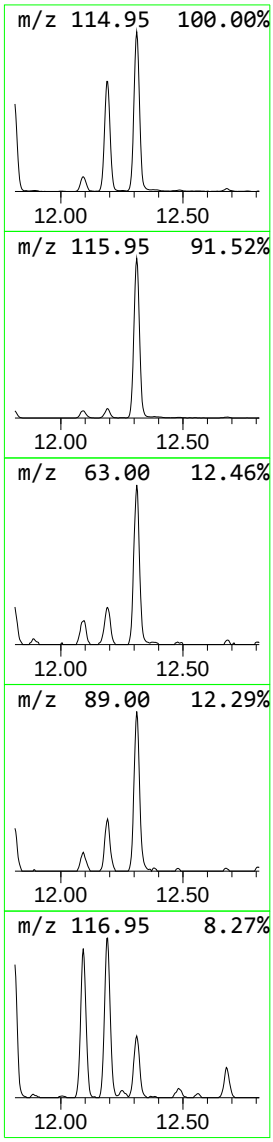
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Indene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.311	25.31 ug/L	462160	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene	116	C9H8	000095-13-6	97
2	Benzene, 1-propynyl-	116	C9H8	000673-32-5	94
3	Benzene, 1-ethynyl-4-methyl-	116	C9H8	000766-97-2	91
4	3-Methylphenylacetylene	116	C9H8	000766-82-5	91
5	1-Propyne, 3-phenyl-	116	C9H8	010147-11-2	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

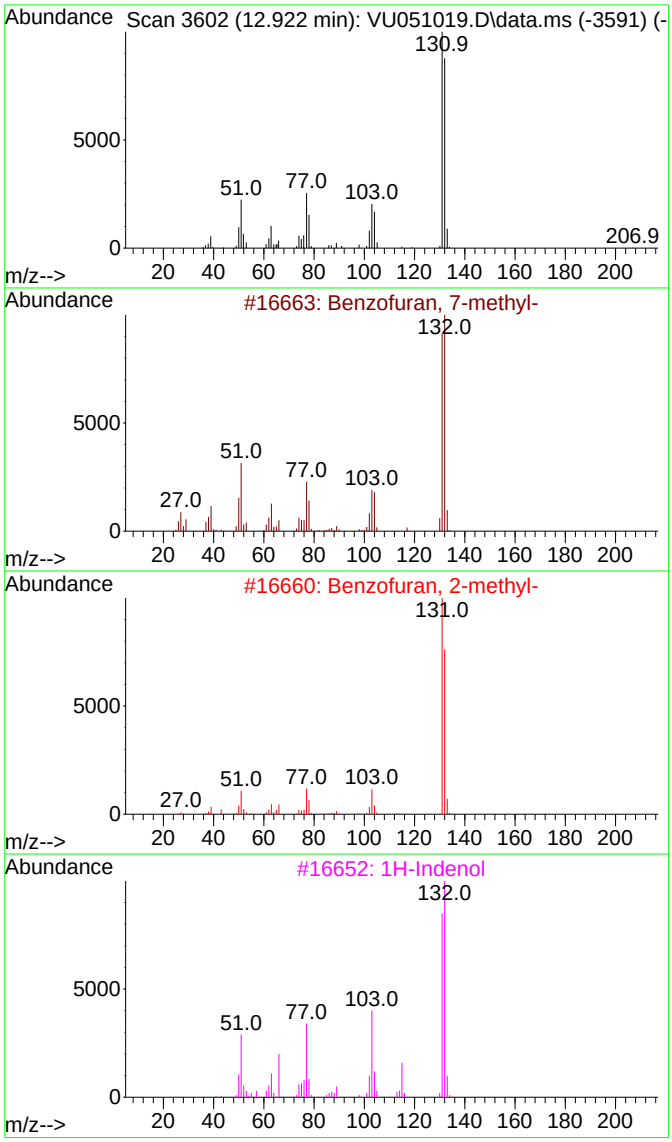
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

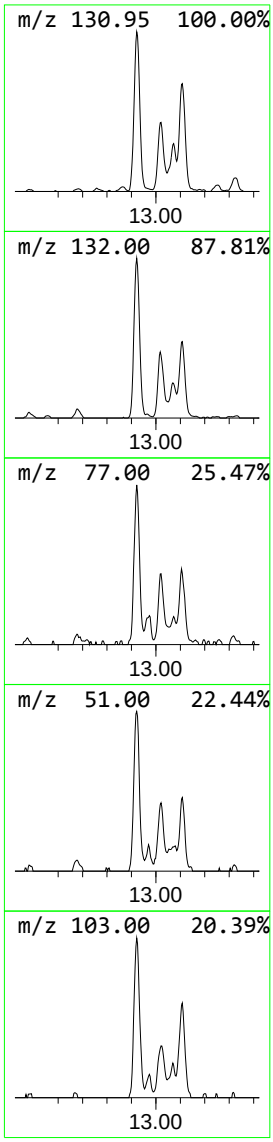
****

Peak Number 4 Benzofuran, 7-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.922	8.50 ug/L	155255	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	1H-Indenol	132	C9H8O	056631-57-3	91
4	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
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ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

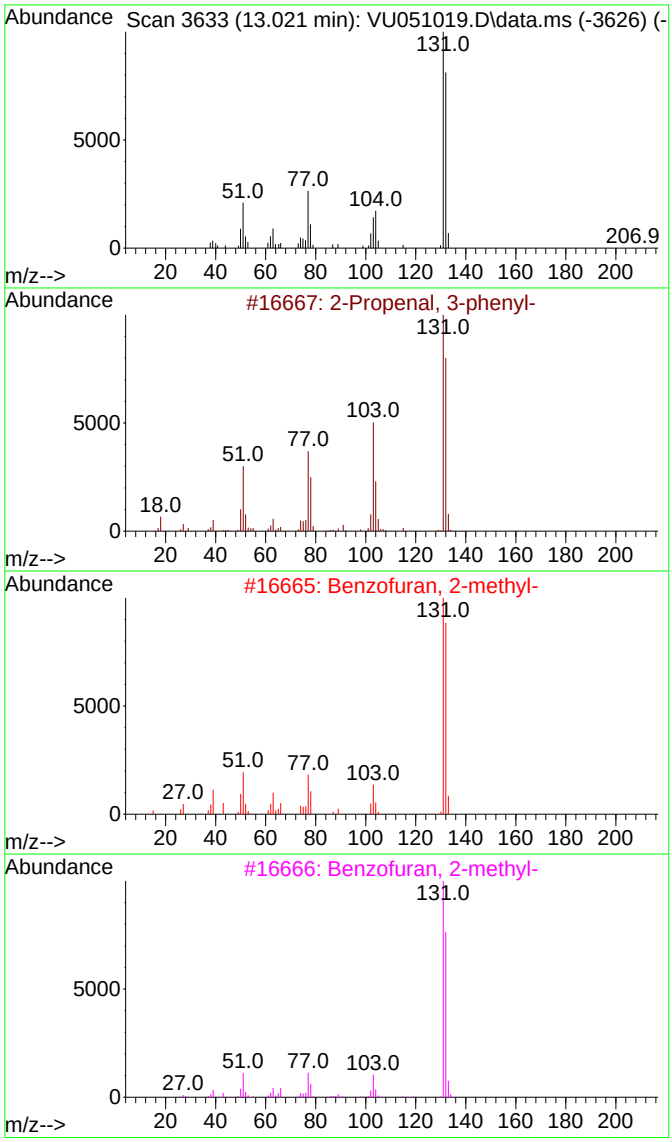
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

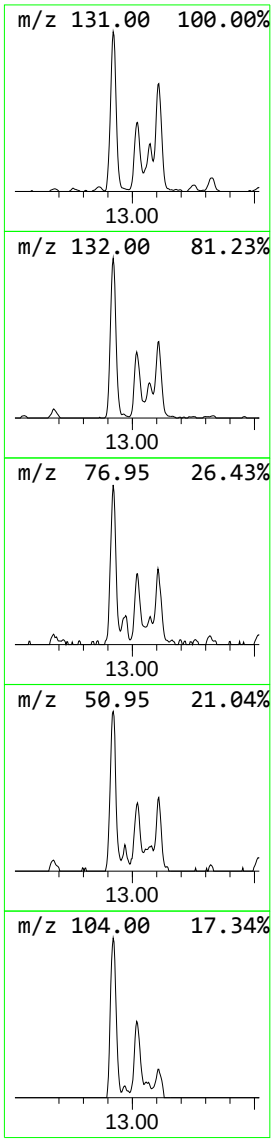
****

Peak Number 5 2-Propenal, 3-phenyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.021	2.88 ug/L	52649	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Propenal, 3-phenyl-	132	C9H8O	000104-55-2	93
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	93
3	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90
5	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

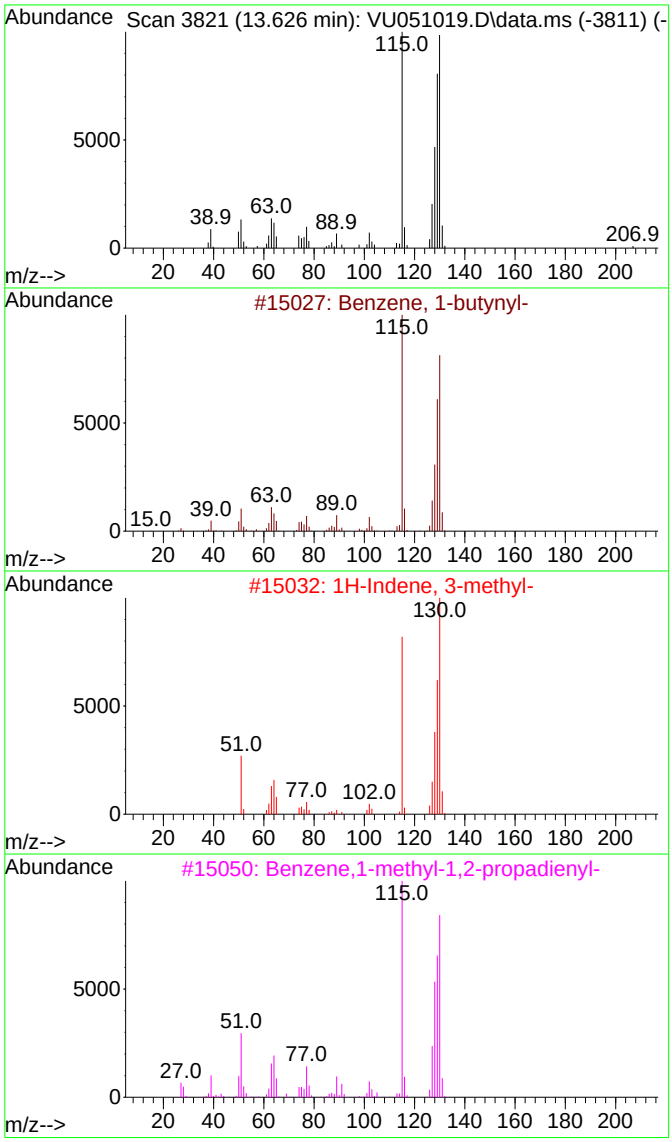
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

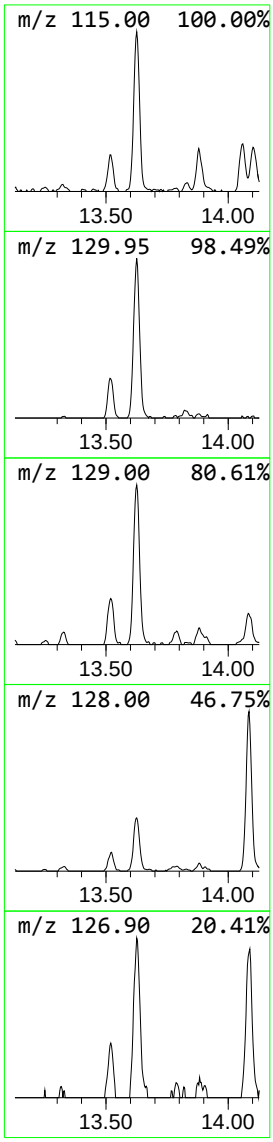
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-butynyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.626	5.71 ug/L	104258	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-butynyl-	130	C10H10	000622-76-4	95
2	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	95
3	Benzene,1-methyl-1,2-propadienyl-	130	C10H10	022433-39-2	94
4	2-Methylindene	130	C10H10	002177-47-1	93
5	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

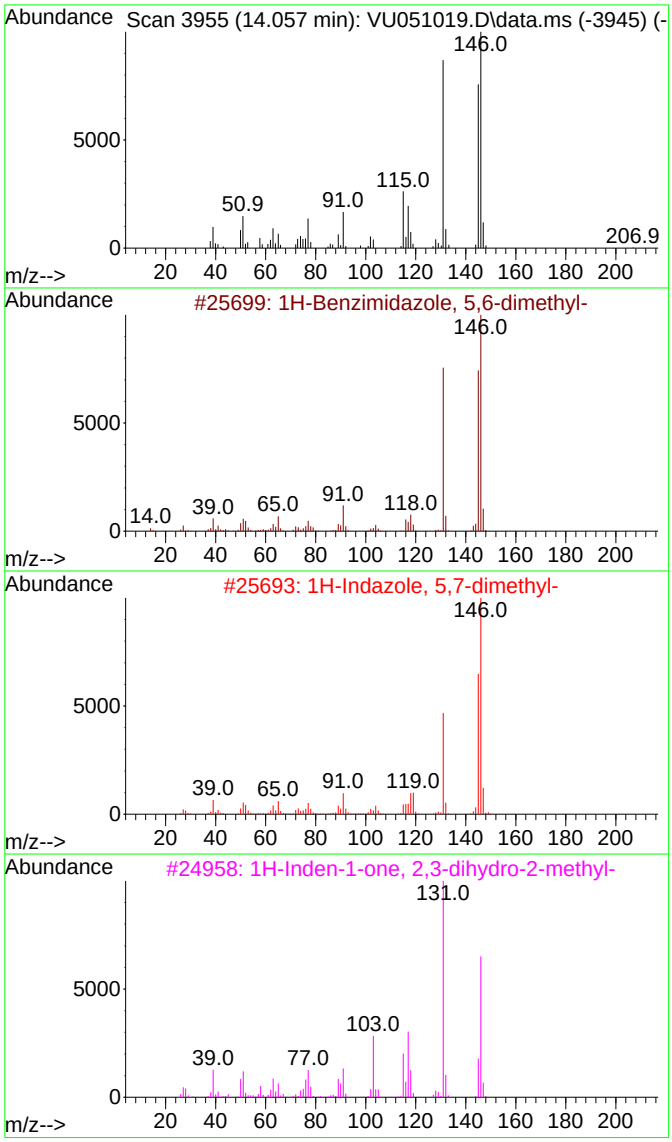
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

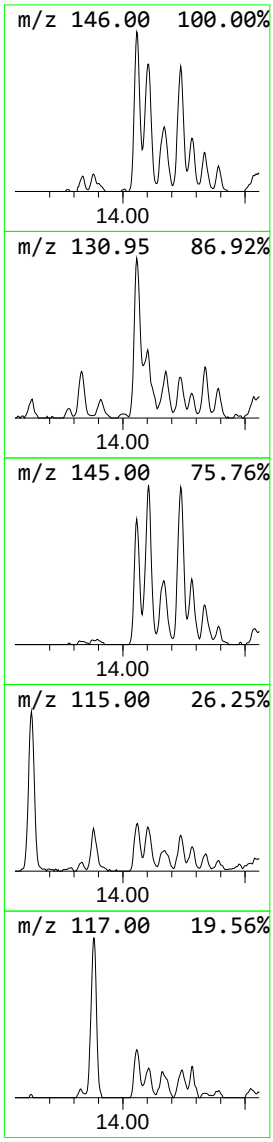
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 1H-Benzimidazole, 5,6-dimet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.057	6.20 ug/L	113195	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	87
2	1H-Indazole, 5,7-dimethyl-	146	C9H10N2	043067-41-0	80
3	1H-Inden-1-one, 2,3-dihydro-2-me...	146	C10H10O	017496-14-9	76
4	2-Amino-4,6-dimethylbenzonitrile	146	C9H10N2	1010461-05-1	64
5	1,2-Dimethylbenzimidazole	146	C9H10N2	002876-08-6	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

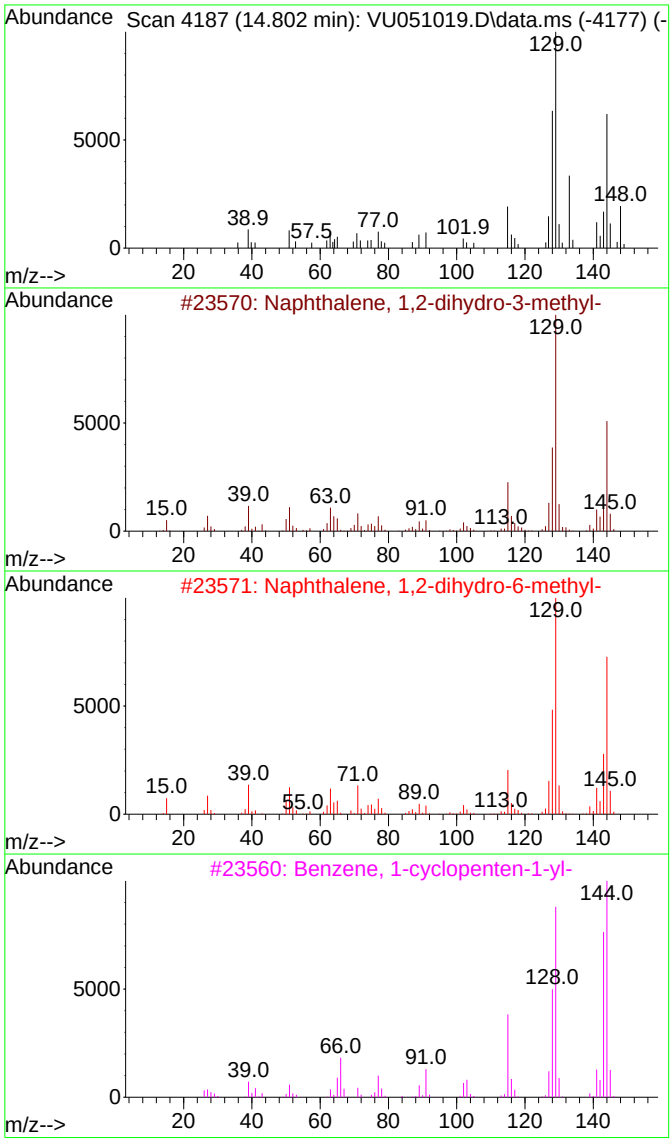
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

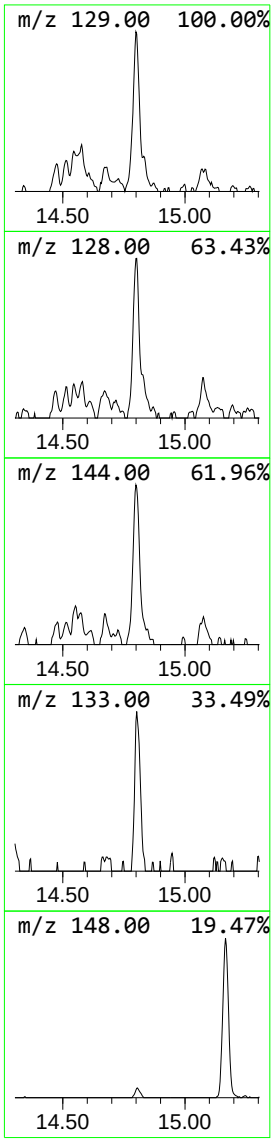
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Naphthalene, 1,2-dihydro-3-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.803	2.60 ug/L	47392	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2-dihydro-3-methyl-	144	C11H12	002717-44-4	93
2	Naphthalene, 1,2-dihydro-6-methyl-	144	C11H12	002717-47-7	92
3	Benzene, 1-cyclopenten-1-yl-	144	C11H12	000825-54-7	81
4	(1-Methylenebut-2-enyl)benzene	144	C11H12	070588-46-4	81
5	1H-Cyclopropa[b]naphthalene, 1a,...	144	C11H12	006571-72-8	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

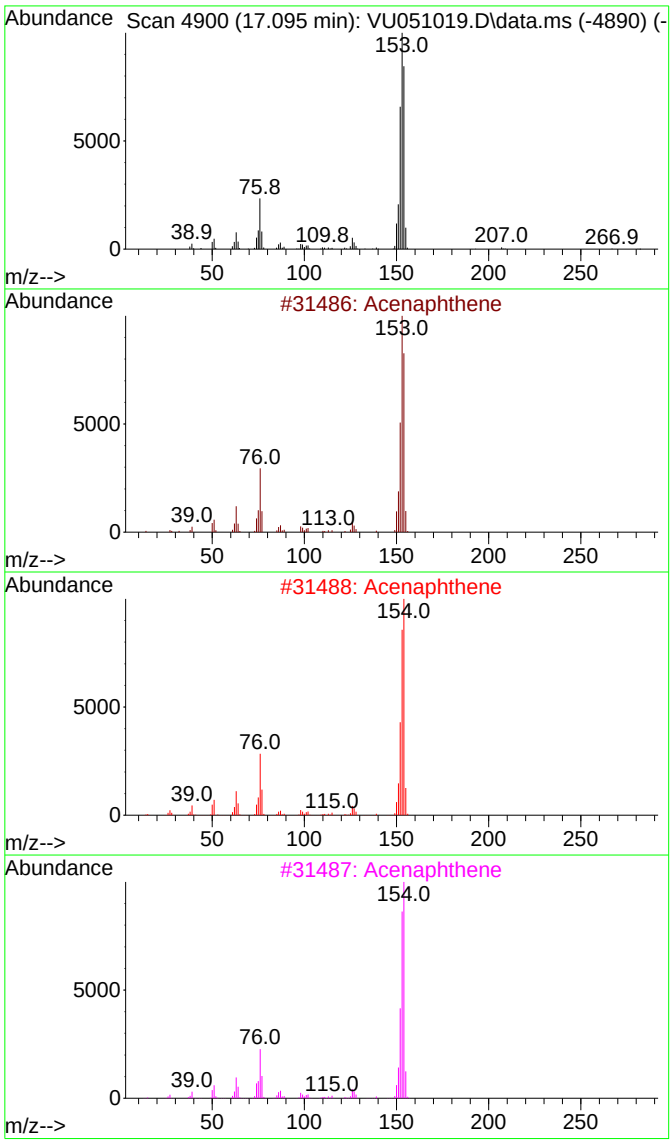
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

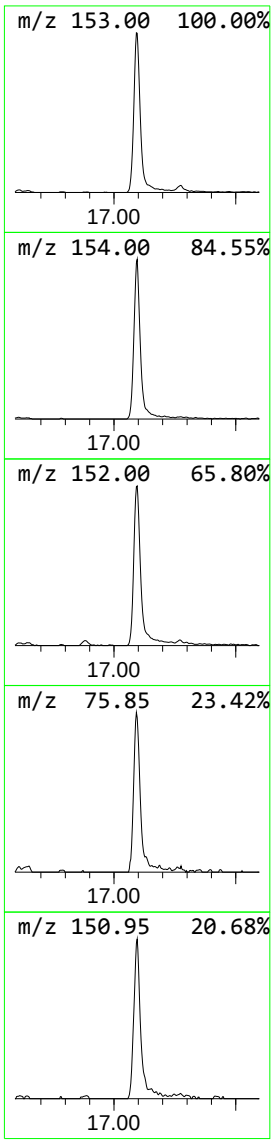
****

Peak Number 9 Hexa-2,4-diyn-1-ylbenzene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.095	20.21 ug/L	368989	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthene	154	C12H10	000083-32-9	70
2	Acenaphthene	154	C12H10	000083-32-9	64
3	Acenaphthene	154	C12H10	000083-32-9	60
4	Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	59
5	3-Fluoro-para-anisaldehyde	154	C8H7FO2	000351-54-2	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU092922\
Data File : VU051019.D
Acq On : 30 Sep 2022 07:24
Operator : JC/MD
Sample : N4859-06
Misc : 5.0mL/MSVOA_U/WATER
ALS Vial : 45 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
E10046

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMULM092922WMA.M
Quant Title : VOC Analysis

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Indane	12.092	5.8	ug/L	106659	3	11.809	913051	50.0
Indene	12.311	25.3	ug/L	462160	3	11.809	913051	50.0
Benzofuran, 7-m...	12.922	8.5	ug/L	155255	3	11.809	913051	50.0
2-Propenal, 3-p...	13.021	2.9	ug/L	52649	3	11.809	913051	50.0
Benzene, 1-buty...	13.626	5.7	ug/L	104258	3	11.809	913051	50.0
1H-Benzimidazol...	14.057	6.2	ug/L	113195	3	11.809	913051	50.0
Naphthalene, 1,...	14.803	2.6	ug/L	47392	3	11.809	913051	50.0
Hexa-2,4-diyn-1...	17.095	20.2	ug/L	368989	3	11.809	913051	50.0