LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

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_V\METHOD\SOMVLM092920WMA.M
Title : VOC Analysis

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.316	63	70	85	rVB	188310	190975	2.90%	0.424%
2	1.579	144	152	164	rBV	161181	181209	2.75%	0.403%
3	2.123	309	321	348	rVB	327243	507305	7.71%	1.128%
4	2.312	377	380	387	rVB3	1035	1171	0.02%	0.003%
5	2.547	429	453	475	rBV	426209	859363	13.05%	1.910%
6	2.708	498	503	508	rBV3	799	840	0.01%	0.002%
7	2.779	508	525	554	rBV	694049	1384881	21.04%	3.078%
8	2.959	573	581	582	rBV5	2434	2857	0.04%	0.006%
9	3.061	596	613	640	rBV	1767455	3506741	53.27%	7.794%
10	3.245	663	670	679	rBV4	4831	9670	0.15%	0.021%
11	3.399	705	718	726	rBV7	4666	10381	0.16%	0.023%
12	3.476	735	742	744	rBV4	2217	2558	0.04%	0.006%
13	3.566	756	770	795	rVB5	27823	91993	1.40%	0.204%
14	3.692	798	809	822	rVV2	18335	41859	0.64%	0.093%
15	3.788	822	839	860	rVV	398522	998418	15.17%	2.219%
16	3.917	868	879	917	rVB	103553	321713	4.89%	0.715%
17	4.377	1004	1022	1048	rBV	237259	592325	9.00%	1.316%
18	4.704	1106	1124	1142	rBV2	35194	89511	1.36%	0.199%
19	4.788	1142	1150	1154	rBV6	1108	1610	0.02%	0.004%
20	4.939	1186	1197	1208	rBV7	2354	5704	0.09%	0.013%
21	5.071	1219	1238	1270	rBV2	604555	1539019	23.38%	3.421%
22	5.521	1370	1378	1392	rVB5	3054	6595	0.10%	0.015%
23	5.640	1402	1415	1441	rBV	514466	1022590	15.53%	2.273%
24	5.933	1494	1506	1522	rVB4	19972	40994	0.62%	0.091%
25	6.093	1541	1556	1580	rBV	338000	687502	10.44%	1.528%
26	6.675	1725	1737	1751	rBV5	4648	9605	0.15%	0.021%
27	6.801	1765	1776	1788	rVB3	5916	11757	0.18%	0.026%
28	6.942	1811	1820	1825	rBV4	2889	4170	0.06%	0.009%
29	7.007	1829	1840	1861	rBV	202450	357622	5.43%	0.795%
30	7.283	1914	1926	1932	rBV3	10320	18974	0.29%	0.042%
31	7.338	1932	1943	1963	rVB	738475	1264628	19.21%	2.811%
32	7.531	2001	2003	2011	rVB5	794	792	0.01%	0.002%
33	7.589	2013	2021	2028	rBV4	4656	6629	0.10%	0.015%
34	7.643	2028	2038	2060	rBV	141254	245465	3.73%	0.546%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

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_V\METHOD\SOMVLM092920WMA.M
Title : VOC Analysis

35	7.807	2084	2089	2098	rVB5	1517	2140	0.03%	0.005%
36	7.897	2110	2117	2125	rBV6	2433	3912	0.06%	0.009%
37	7.997	2142	2148	2149	rBV4	844	793	0.01%	0.002%
38	8.003	2149	2150	2157	rVB4	1176	938	0.01%	0.002%
39	8.109	2173	2183	2216	rBV	369056	667127	10.13%	1.483%
40	8.524	2305	2312	2315	rVB5	734	803	0.01%	0.002%
41	8.589	2326	2332	2335	rBV5	1384	1526	0.02%	0.003%
42	8.688	2356	2363	2371	rVB5	2732	4584	0.07%	0.010%
43	8.810	2394	2401	2409	rBV6	1584	2456	0.04%	0.005%
44	8.872	2409	2420	2444	rBV	769553	1259890	19.14%	2.800%
45	9.035	2461	2471	2486	rBV	47589	76791	1.17%	0.171%
46	9.158	2499	2509	2523	rBV	227894	374056	5.68%	0.831%
47	9.225	2525	2530	2545	rVB5	10665	19748	0.30%	0.044%
48	9.495	2603	2614	2621	rBV6	5089	8227	0.12%	0.018%
49	9.566	2625	2636	2654	rBV	58271	97493	1.48%	0.217%
50	9.701	2671	2678	2681	rBV5	1697	2372	0.04%	0.005%
51	9.952	2742	2756	2774	rBV	171019	274392	4.17%	0.610%
52	10.039	2775	2783	2791	rBV6	5130	7995	0.12%	0.018%
53	10.145	2809	2816	2828	rVB2	21650	36327	0.55%	0.081%
54	10.238	2833	2845	2864	rVV	452037	713536	10.84%	1.586%
55	10.373	2876	2887	2903	rBV	851332	1267443	19.25%	2.817%
56	10.470	2905	2917	2935	rBV	2998146	6296827	95.65%	13.995%
57	10.560	2935	2945	2955	rBV	1562716	2254994	34.25%	5.012%
58	10.759	2995	3007	3026	rVB	1032741	1558240	23.67%	3.463%
59	10.936	3044	3062	3087	rVB	4433853	6583269	100.00%	14.632%
60	11.077	3093	3106	3111	rBV2	29194	53645	0.81%	0.119%
61	11.167	3127	3134	3141	rBV5	16797	26071	0.40%	0.058%
62	11.215	3141	3149	3156	rVV	57547	84359	1.28%	0.187%
63	11.270	3156	3166	3182	rVV	907319	1368635	20.79%	3.042%
64	11.357	3182	3193	3216	rVB	910835	1411891	21.45%	3.138%
65	11.482	3228	3232	3241	rVB5	14311	18122	0.28%	0.040%
66	11.550	3242	3253	3262	rBV3	232886	485808	7.38%	1.080%
67	11.646	3274	3283	3304	rVB4	1031379	1983867	30.13%	4.409%
68	11.769	3314	3321	3326	rBV2	15196	21212	0.32%	0.047%
69	11.810	3327	3334	3338	rBV	61049	79826	1.21%	0.177%
70	11.842	3339	3344	3362	rVB	91135	138253	2.10%	0.307%
71	11.932	3362	3372	3376	rBV	201845	286691	4.35%	0.637%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

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_V\METHOD\SOMVLM092920WMA.M
Title : VOC Analysis

72	12.039	3395	3405	3426	rVB	380970	588769	8.94%	1.309%
73	12.141	3429	3437	3440	rBV	25714	37797	0.57%	0.084%
74	12.280	3468	3480	3488	rBV8	15685	30977	0.47%	0.069%
75	12.338	3488	3498	3508	rVB	88013	130513	1.98%	0.290%
76	12.434	3508	3528	3536	rBV	244033	381361	5.79%	0.848%
77	12.486	3536	3544	3562	rVB	368365	557984	8.48%	1.240%
78	12.572	3563	3571	3576	rBV	24207	34593	0.53%	0.077%
79	12.608	3577	3582	3586	rBV4	20283	23039	0.35%	0.051%
80	12.746	3616	3625	3638	rVB	115089	174153	2.65%	0.387%
81	12.817	3638	3647	3655	rBV4	18299	27527	0.42%	0.061%
82	12.903	3655	3674	3703	rBV2	278832	546760	8.31%	1.215%
83	13.022	3703	3711	3718	rBV	38909	54172	0.82%	0.120%
84	13.186	3755	3762	3770	rBV3	10133	14748	0.22%	0.033%
85	13.231	3773	3776	3783	rVB5	3390	3715	0.06%	0.008%
86	13.286	3783	3793	3802	rBV2	178483	302923	4.60%	0.673%
87	13.421	3828	3835	3847	rVB	28418	40487	0.61%	0.090%
88	13.527	3858	3868	3879	rBV	220740	333282	5.06%	0.741%
89	13.768	3924	3943	3956	rBV4	46280	98241	1.49%	0.218%
90	13.929	3984	3993	4007	rVB3	25030	43756	0.66%	0.097%
91	14.112	4042	4050	4064	rBV7	11393	25342	0.38%	0.056%
92	14.257	4087	4095	4104	rBV3	9408	14911	0.23%	0.033%
93	14.662	4208	4221	4232	rBV3	13736	25271	0.38%	0.056%
94	14.791	4256	4261	4263	rBV3	1210	1203	0.02%	0.003%
95	14.849	4274	4279	4282	rBV3	3043	3695	0.06%	0.008%
96	14.968	4310	4316	4317	rBV4	1146	1001	0.02%	0.002%
97	15.067	4345	4347	4351	rVV3	1052	893	0.01%	0.002%
98	15.315	4420	4424	4429	rBV8	1248	1241	0.02%	0.003%
99	15.344	4430	4433	4436	rBV5	1354	1143	0.02%	0.003%
100	15.440	4460	4463	4470	rVB5	1302	1243	0.02%	0.003%

Sum of corrected areas: 44992425

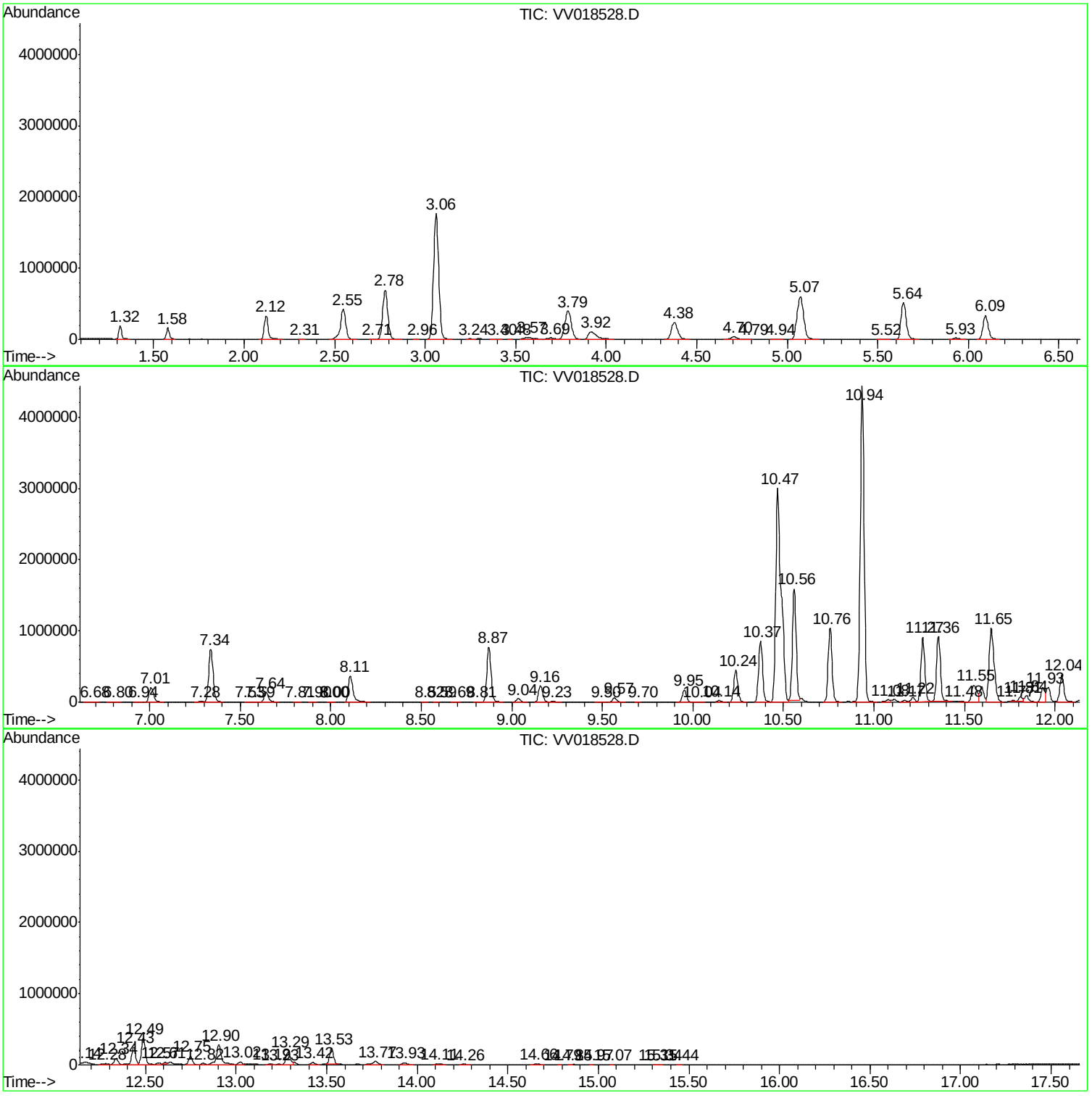
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

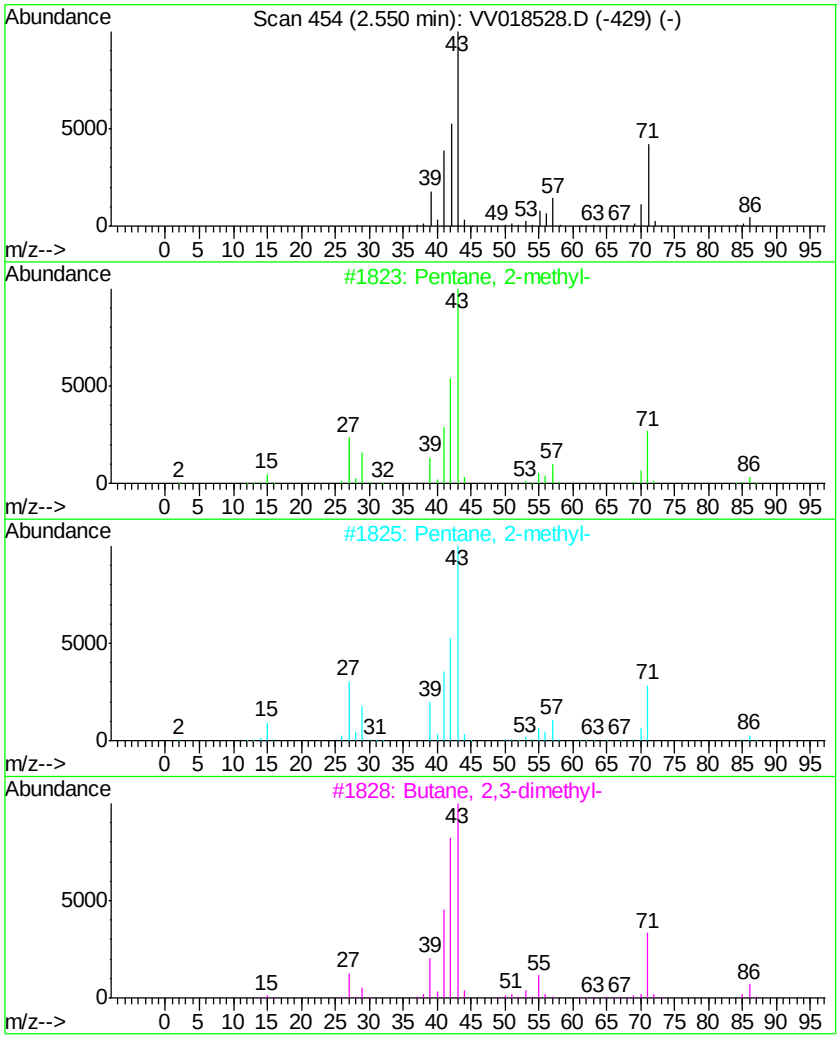
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

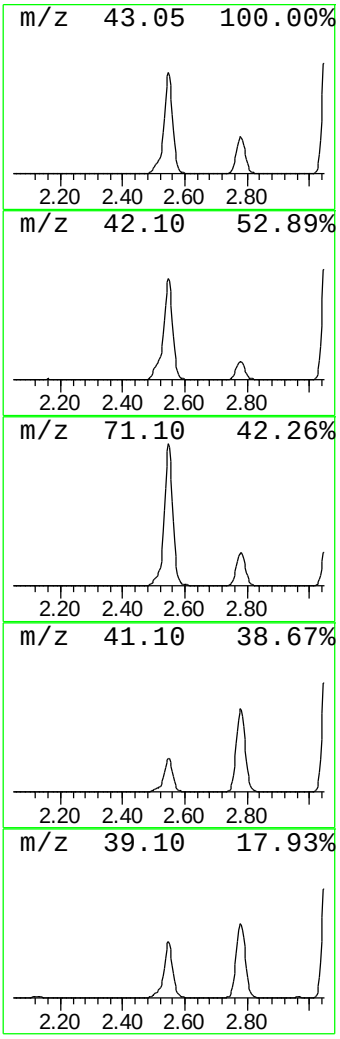
****

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.55	42.02 ug/L	859363	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
2	Pentane, 2-methyl-	86	C6H14	000107-83-5	91
3	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	72
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	49
5	1-Butanol, 2,3-dimethyl-	102	C6H14O	019550-30-2	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

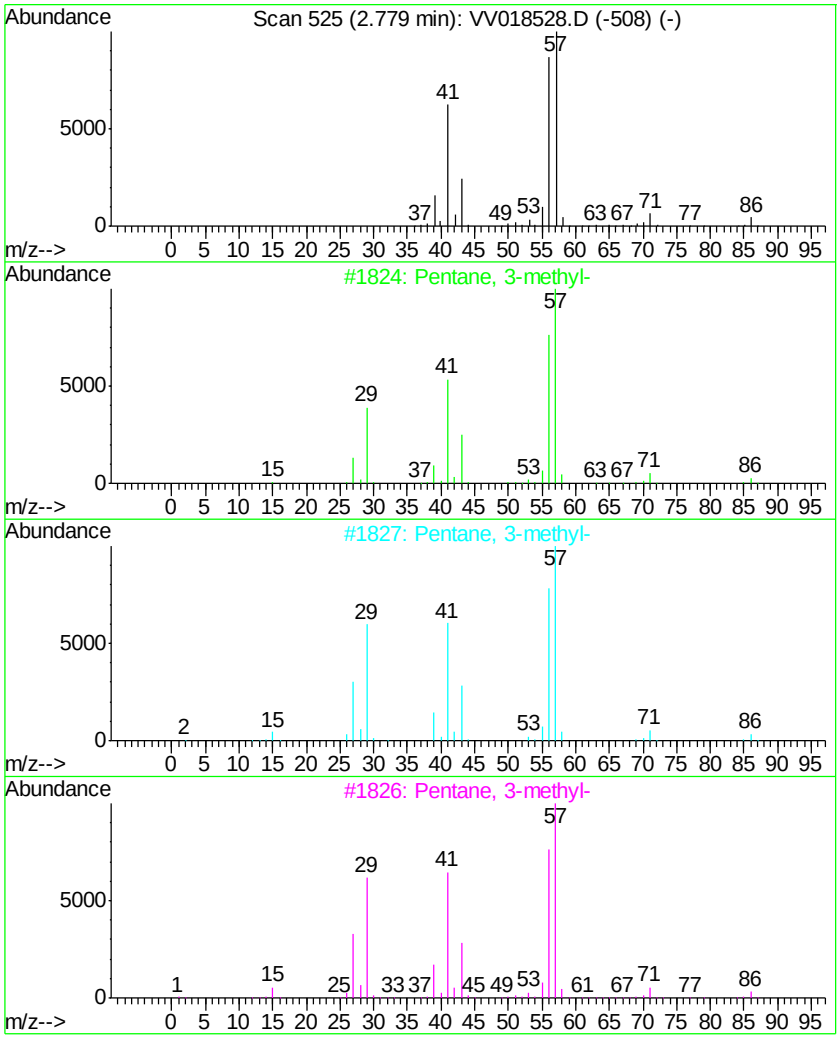
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

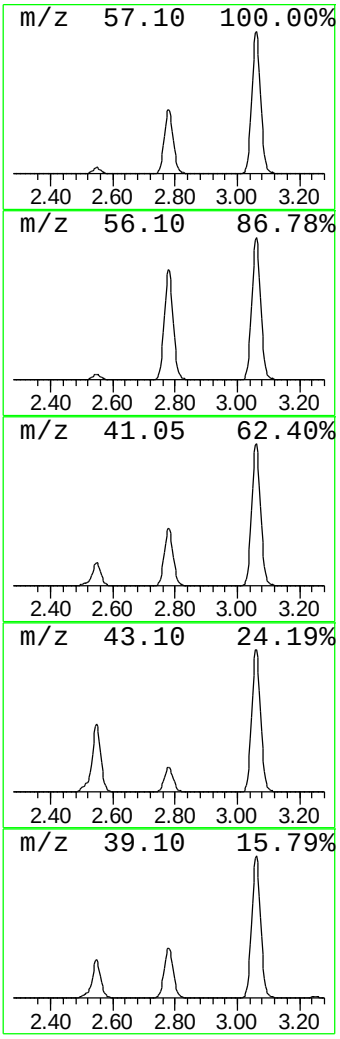
****

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.78	67.71 ug/L	1384880	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
4	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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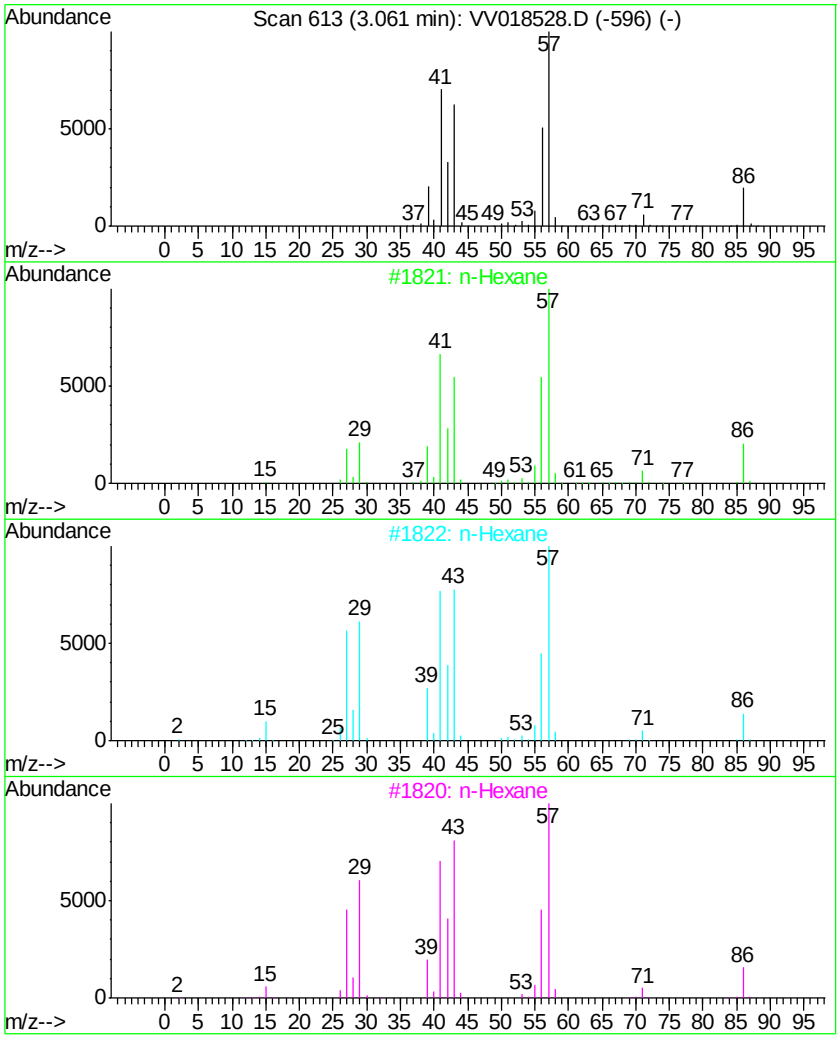
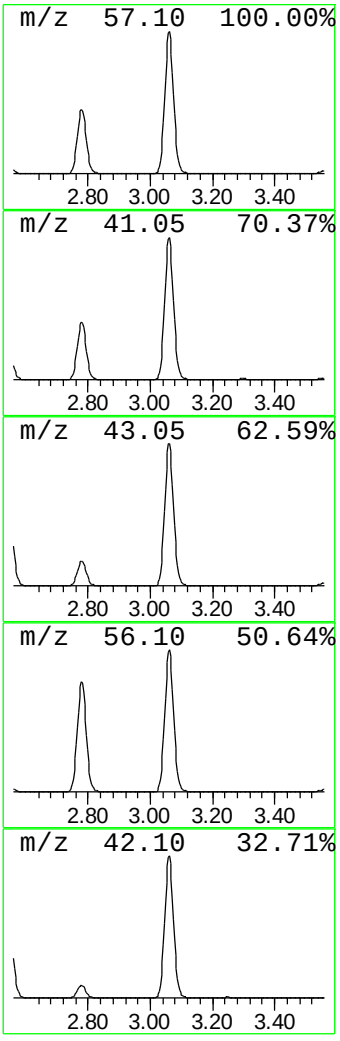
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.06	171.46 ug/L	3506740	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	94
2	n-Hexane	86	C6H14	000110-54-3	78
3	n-Hexane	86	C6H14	000110-54-3	78
4	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	40
5	Butanal, 2-methyl-	86	C5H10O	000096-17-3	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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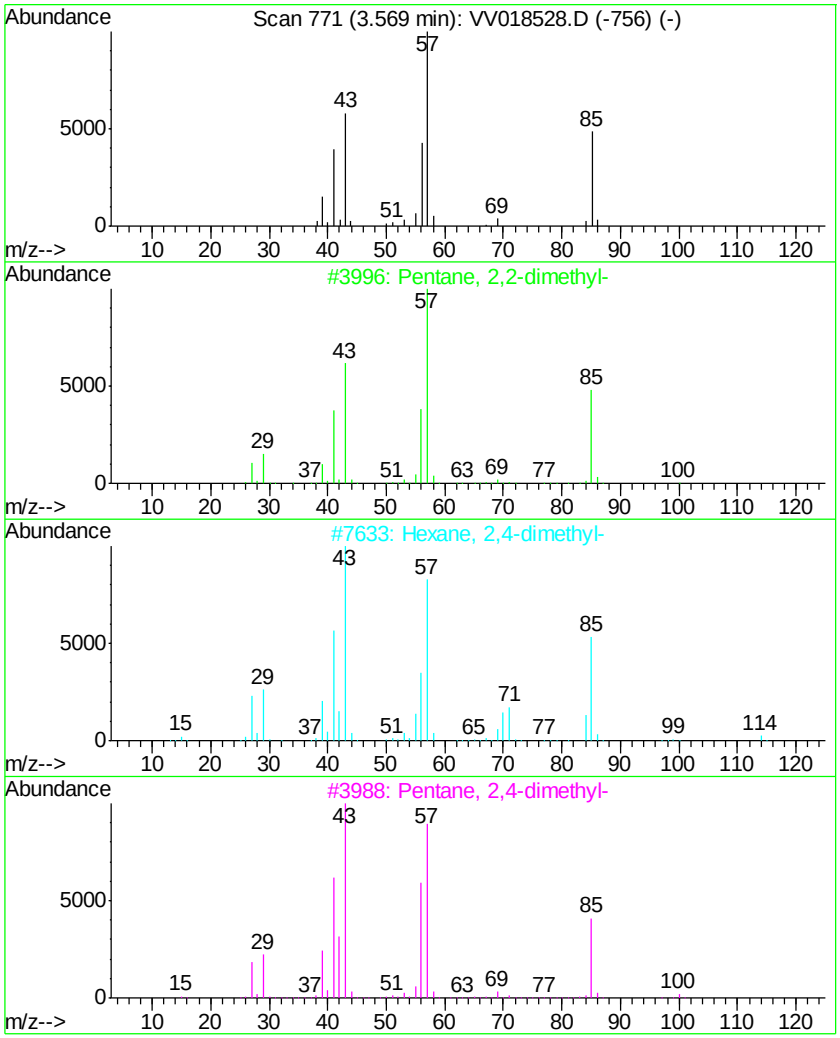
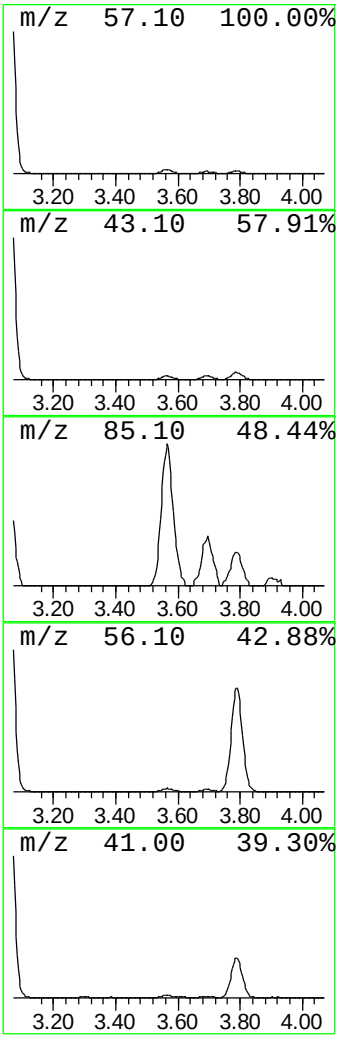
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.57	4.50 ug/L	91993	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
2	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	78
3	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
4	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64
5	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

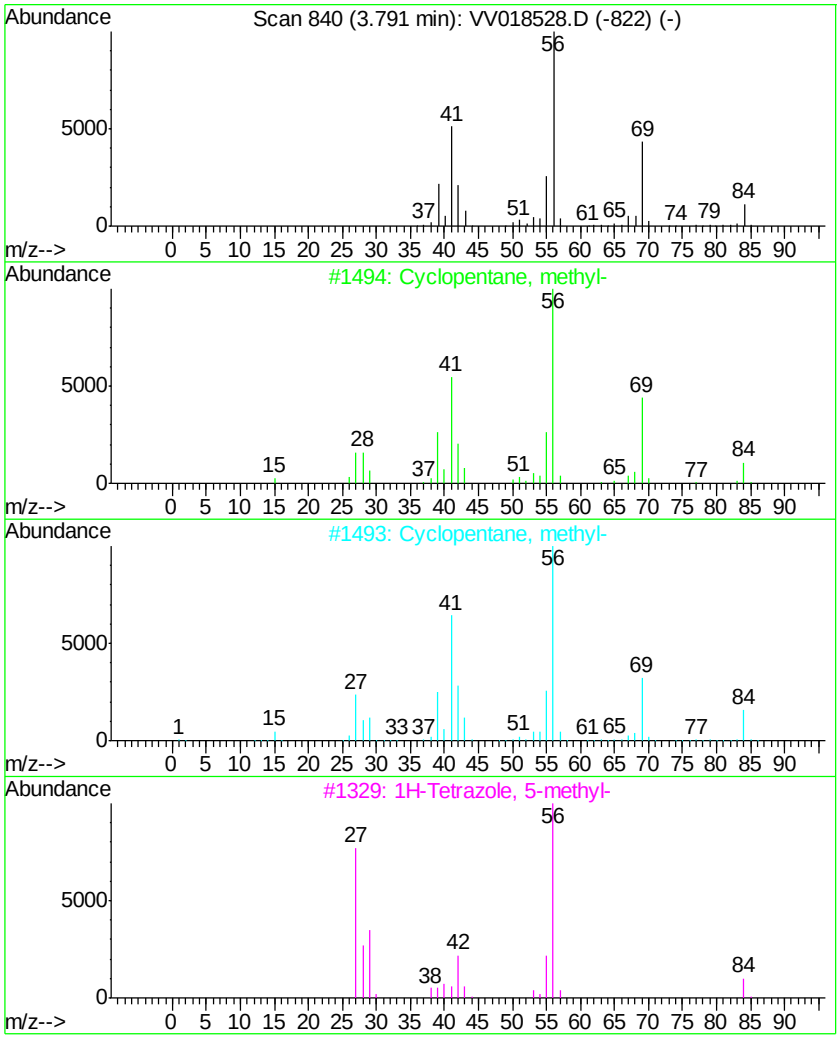
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

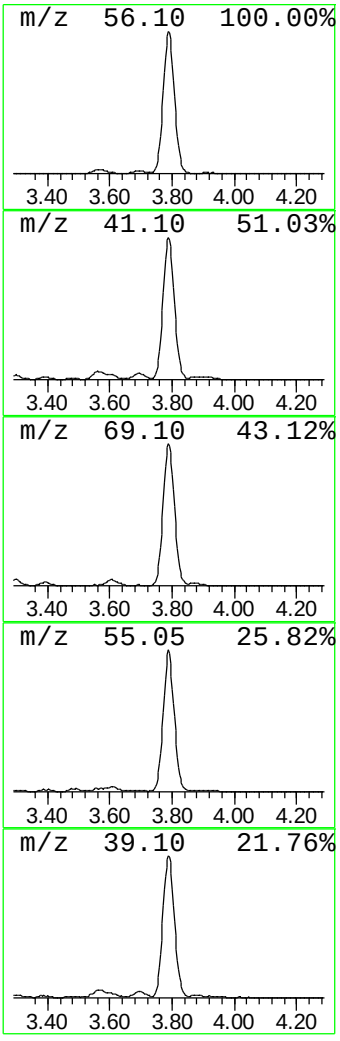
****

Peak Number 5 (DEL) Alkane: Cyclic3.79 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.79	48.82 ug/L	998418	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
4	Cyclohexane	84	C6H12	000110-82-7	78
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

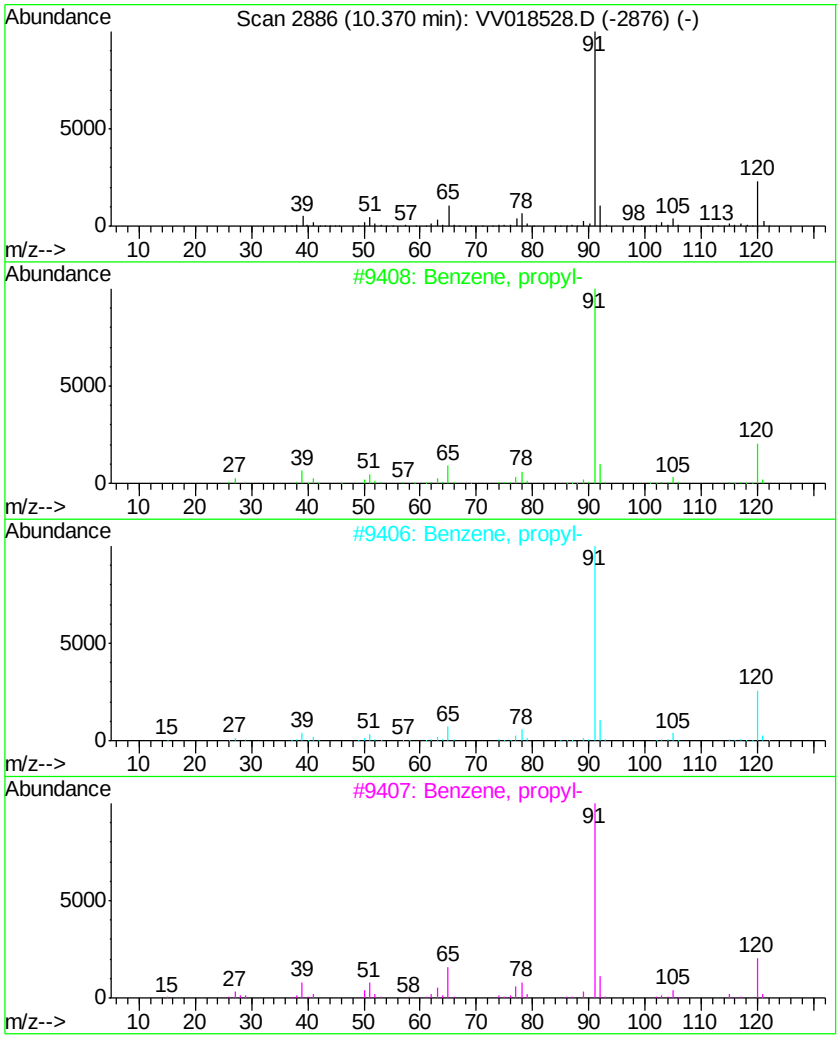
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

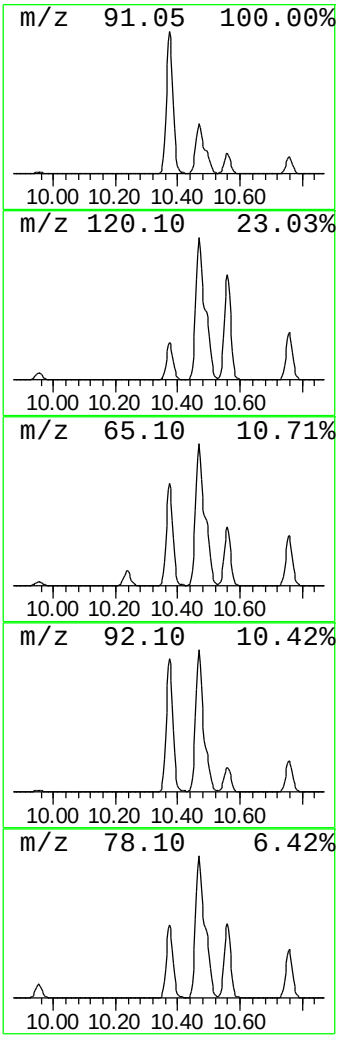
****

Peak Number 7 Benzene, propyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.37	46.30 ug/L	1267440	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	90
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	86
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	74
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

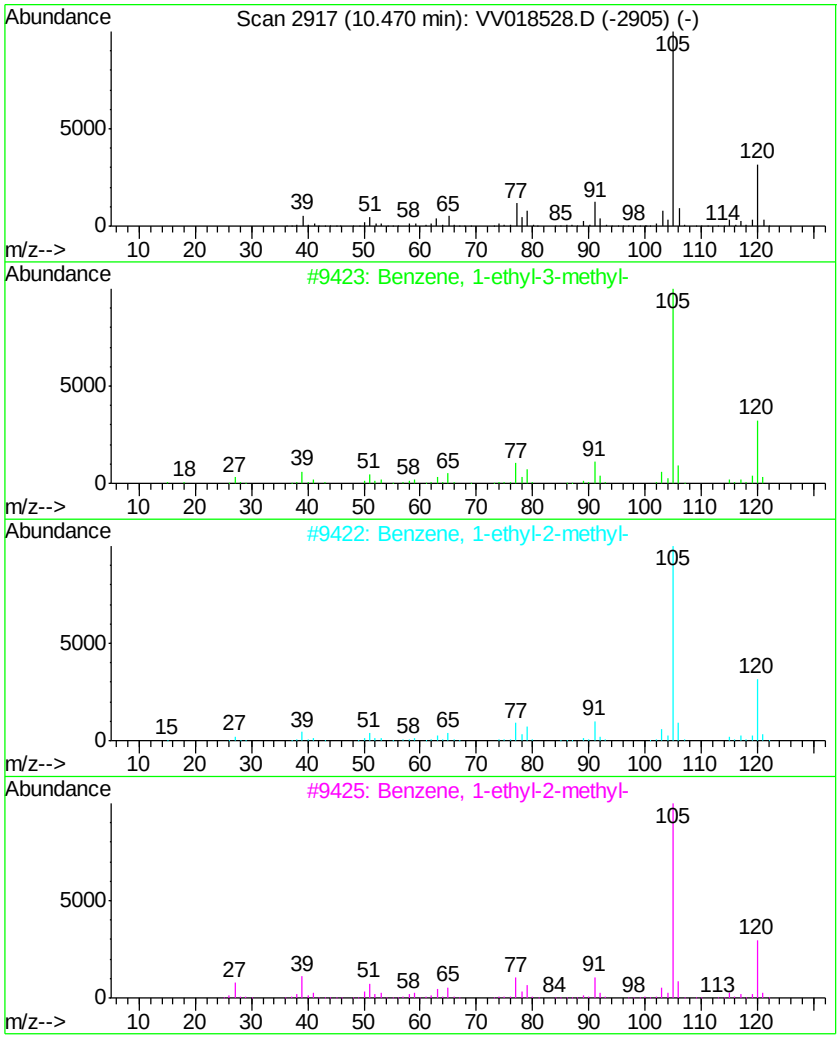
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

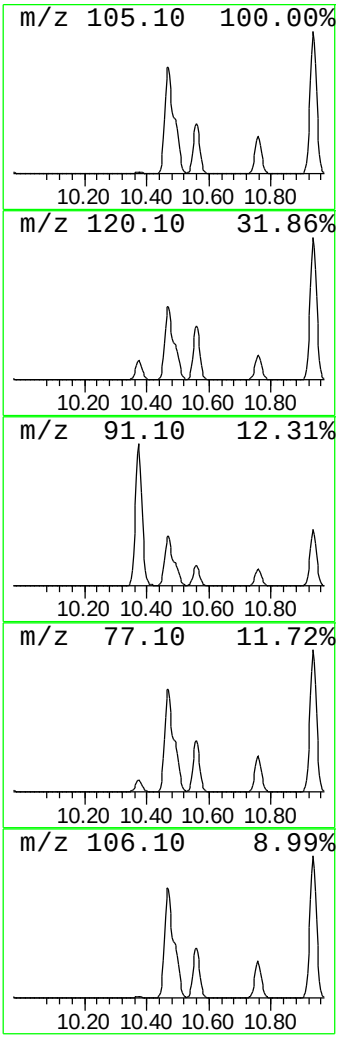
****

Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.47	230.04 ug/L	6296830	1,4-Dichlorobenzene-d4	11.27

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-ethyl-2-methyl-	120	C9H12	000611-14-3	95
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	95
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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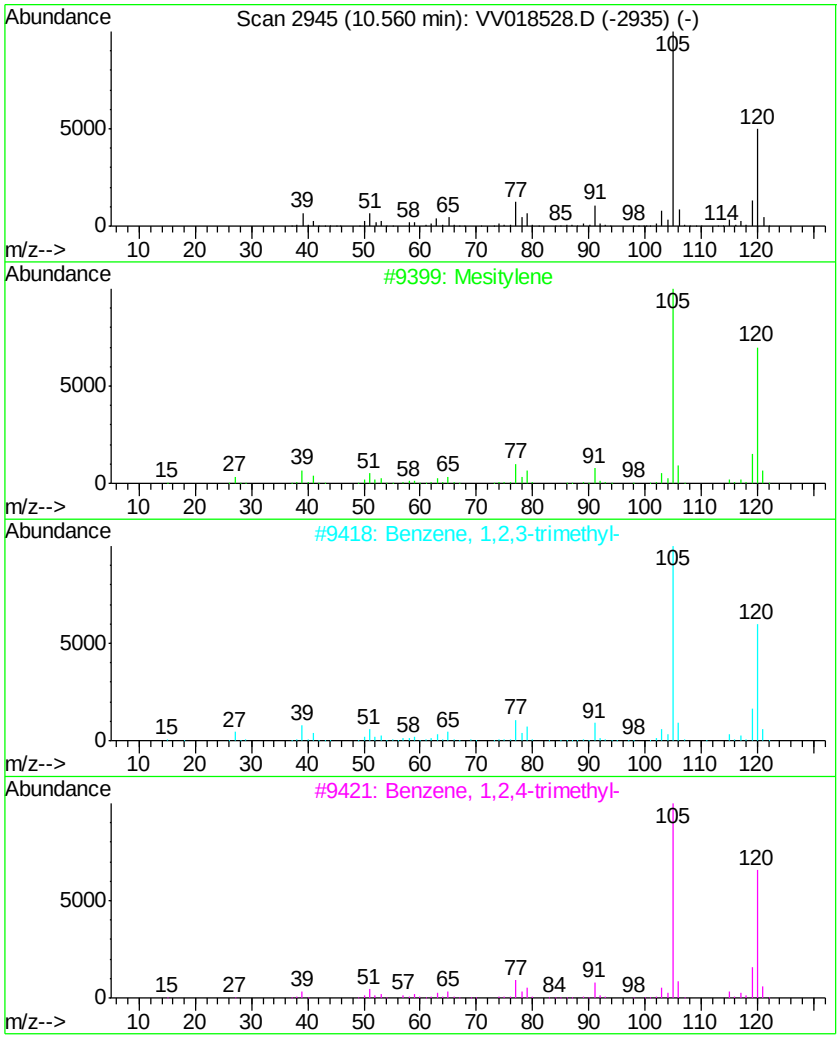
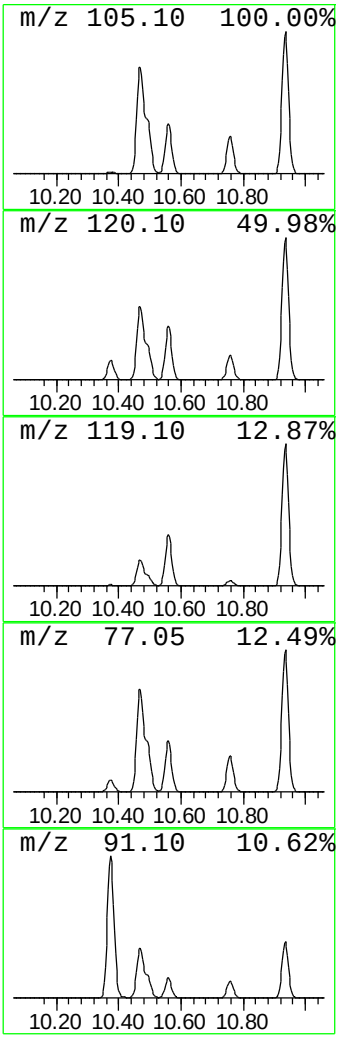
Peak Number 9 Mesitylene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.56	82.38 ug/L	2254990	1,4-Dichlorobenzene-d4	11.27

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	97
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
5	Mesitylene	120	C9H12	000108-67-8	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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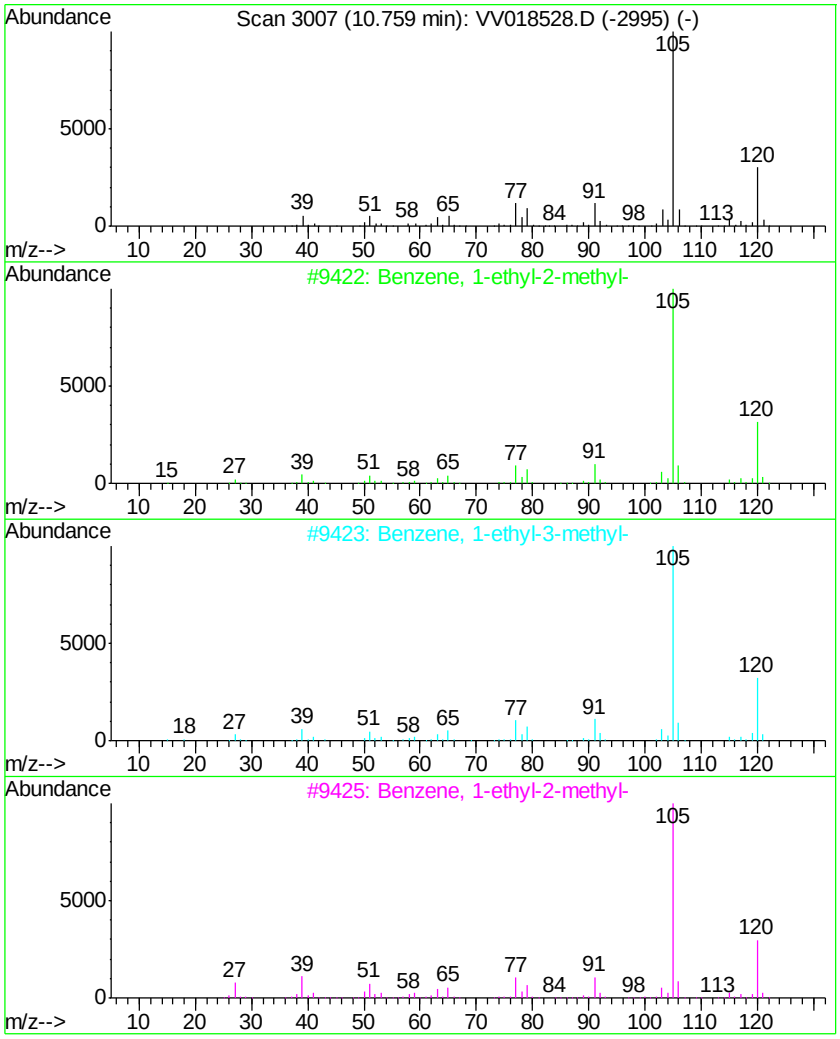
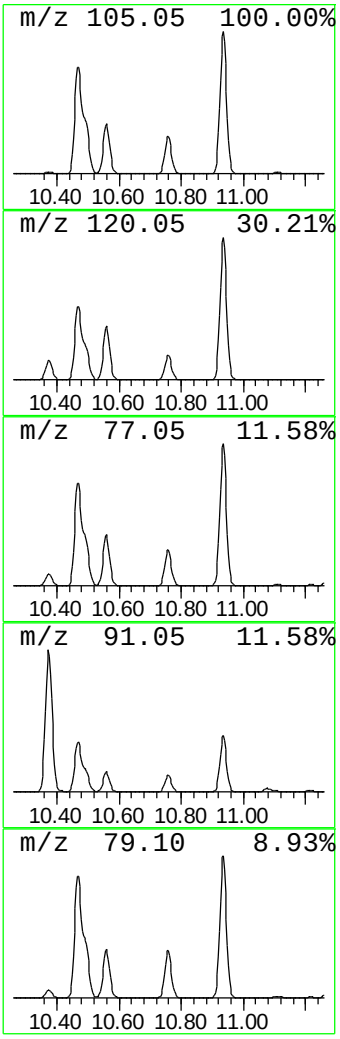
Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.76	56.93 ug/L	1558240	1,4-Dichlorobenzene-d4	11.27

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-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
5	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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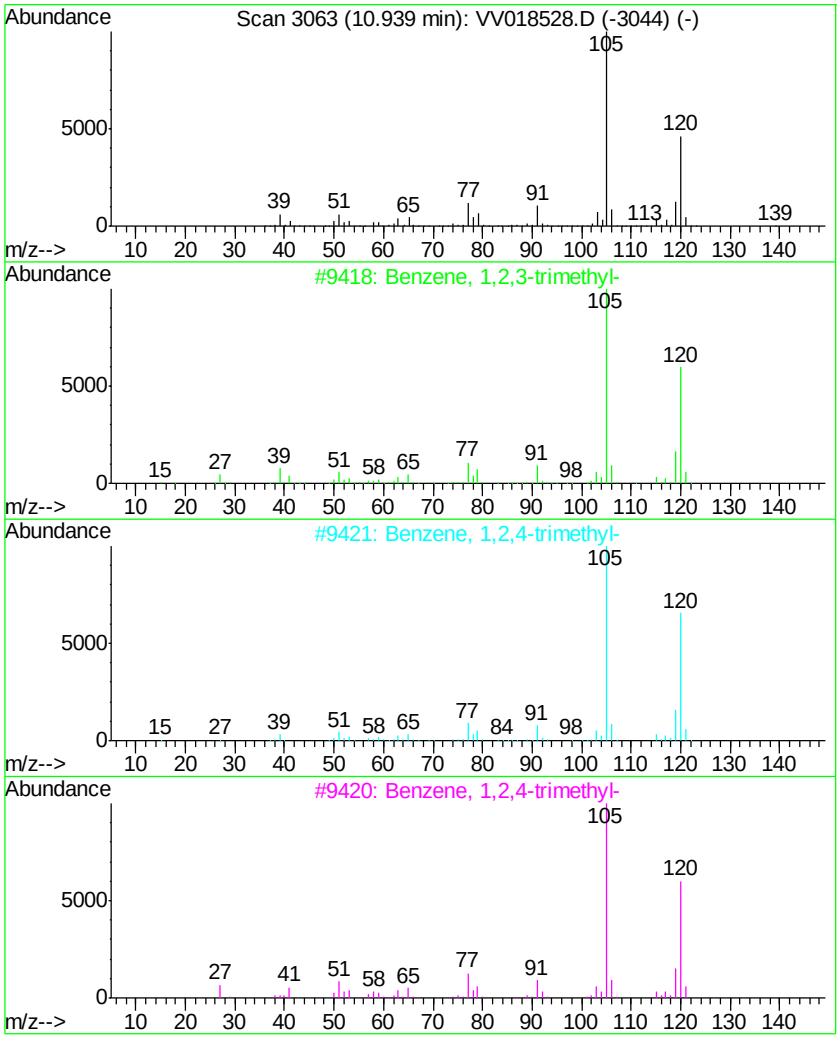
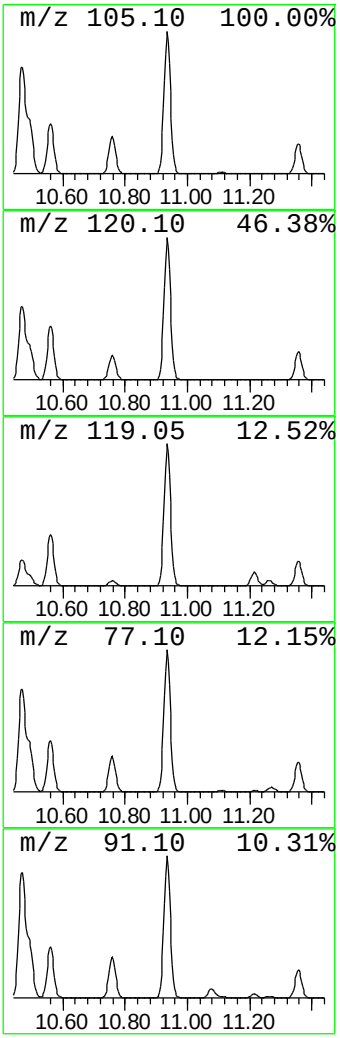
Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.94	240.51 ug/L	6583270	1,4-Dichlorobenzene-d4	11.27

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





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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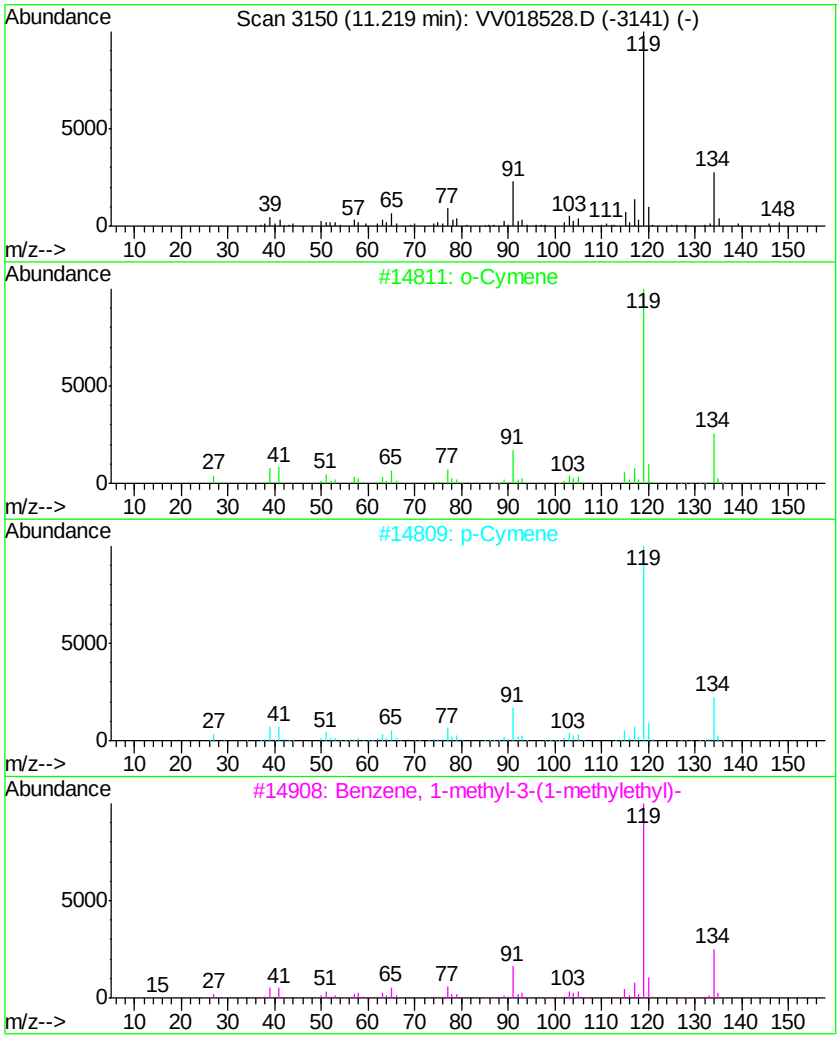
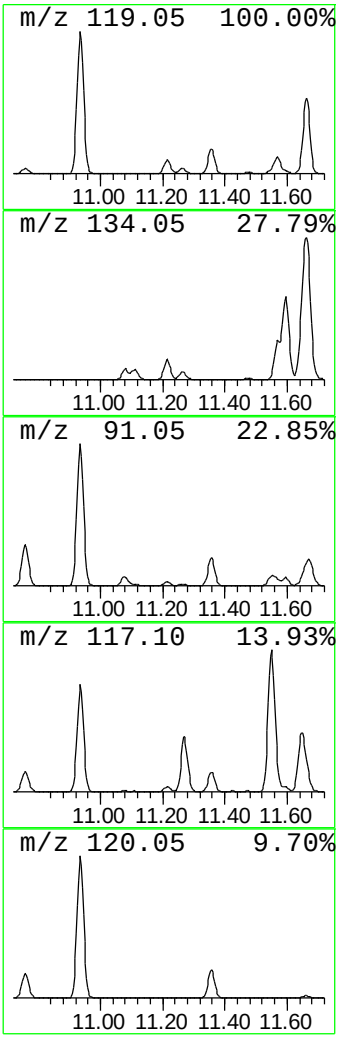
Peak Number 12 o-Cymene Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.22	3.08 ug/L	84359	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	p-Cymene	134	C10H14	000099-87-6	97
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	97
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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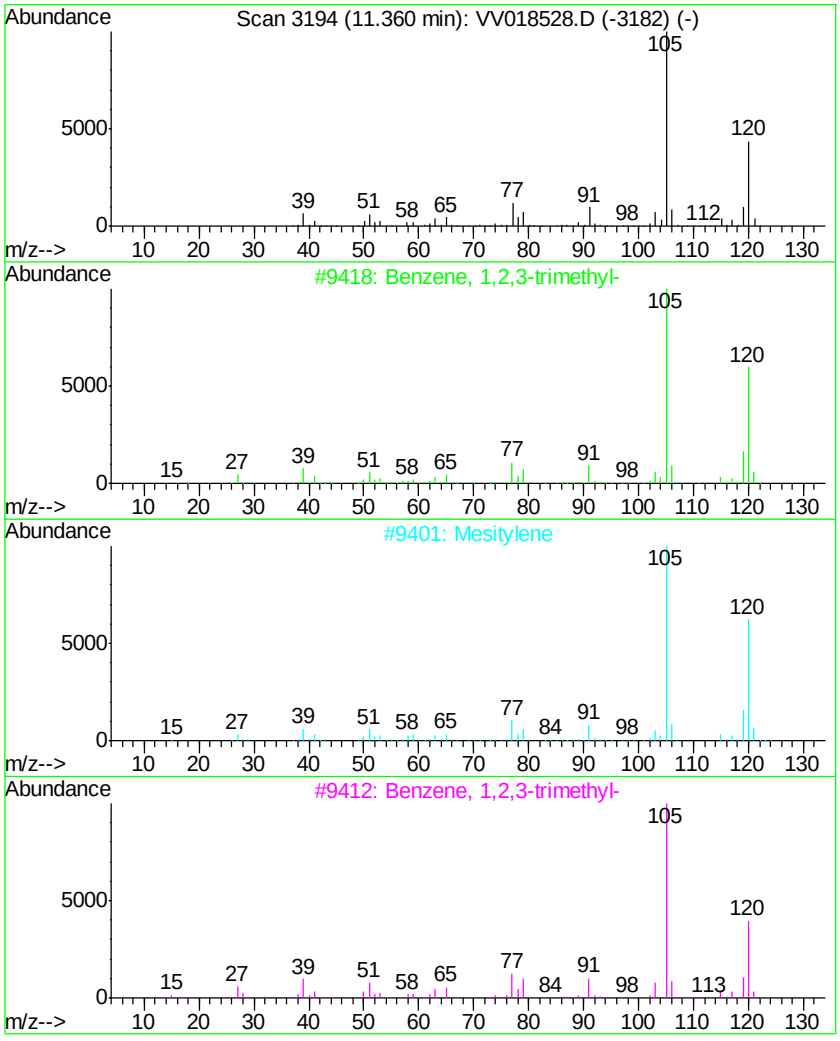
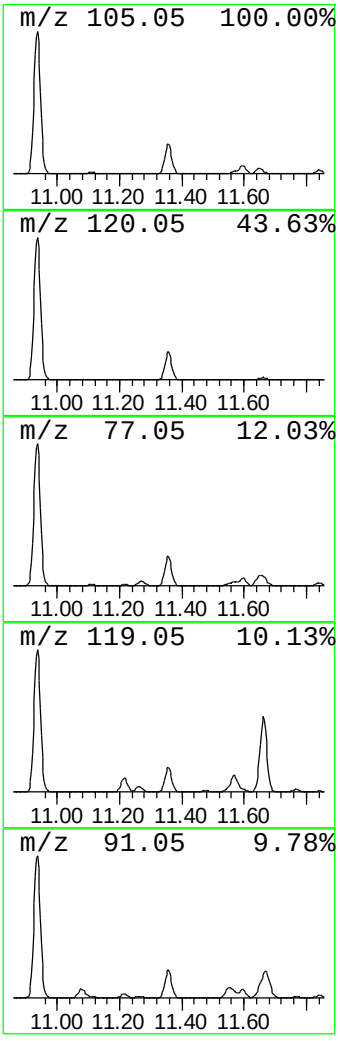
Peak Number 13 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.36	51.58 ug/L	1411890	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Mesitylene	120	C9H12	000108-67-8	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4	Mesitylene	120	C9H12	000108-67-8	93
5	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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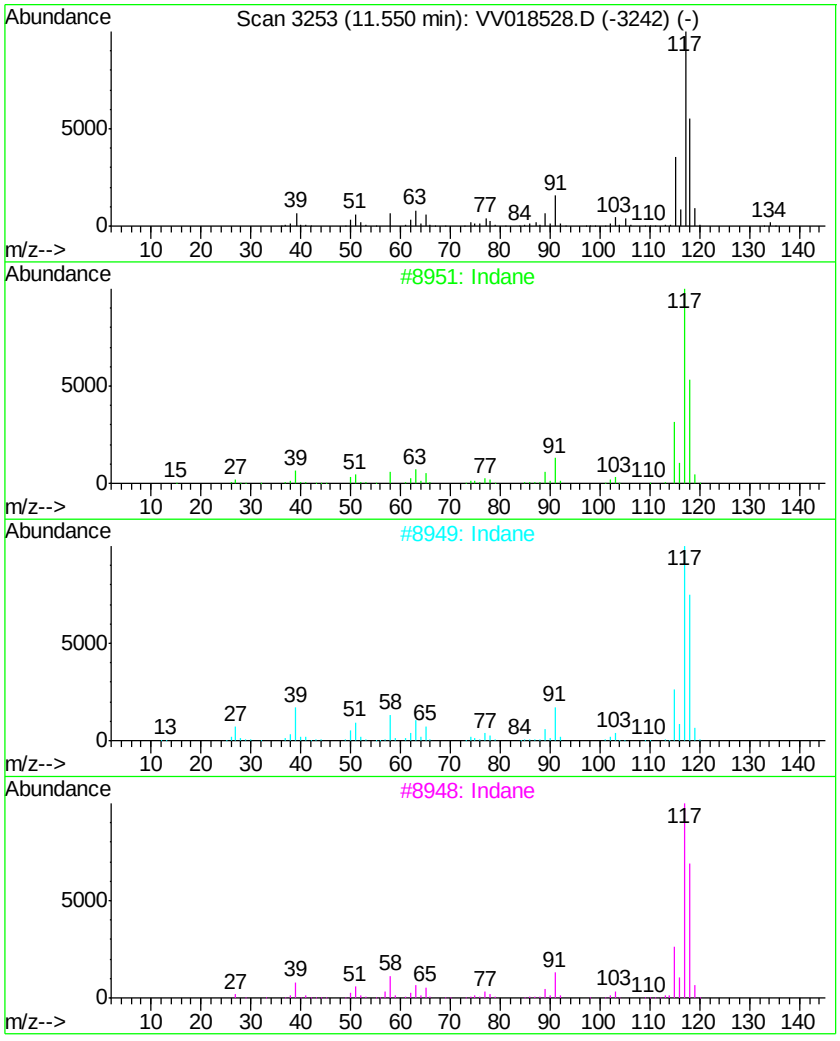
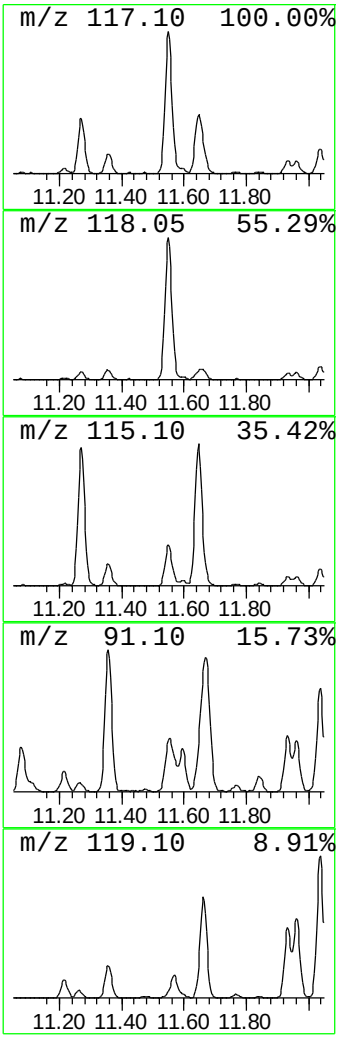
Peak Number 14 Indane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.55	17.75 ug/L	485808	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	94
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
5	Benzene, 2-propenyl-	118	C9H10	000300-57-2	74





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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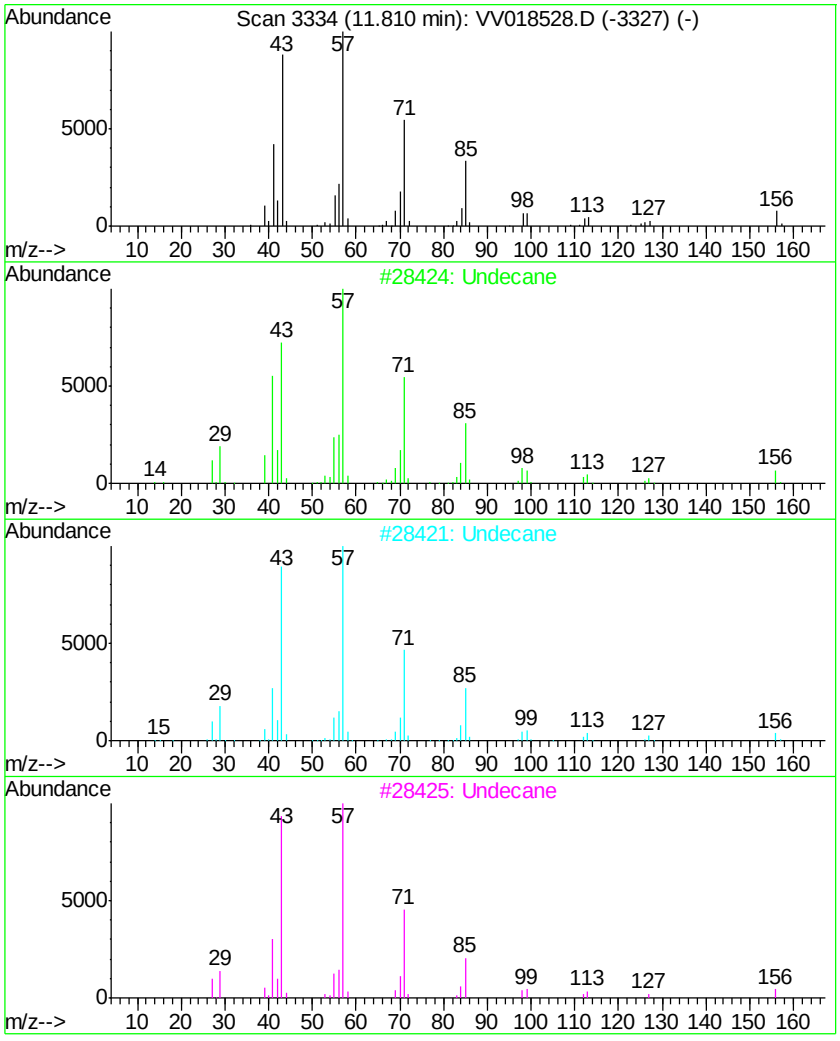
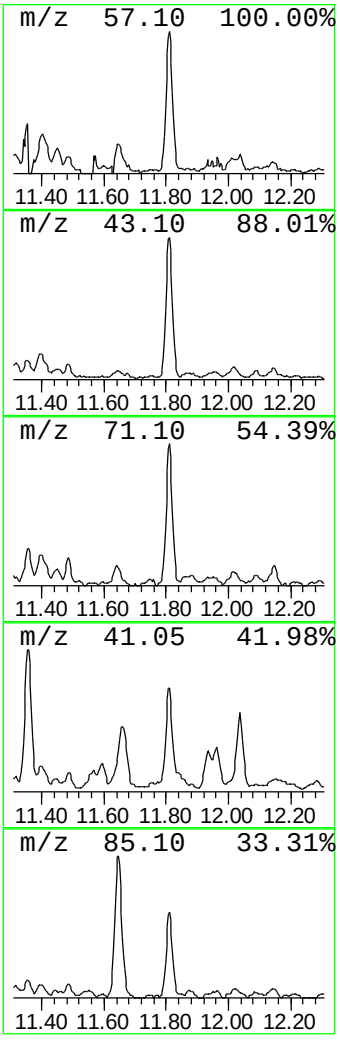
Peak Number 15 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	2.92 ug/L	79826	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	96
2	Undecane	156	C11H24	001120-21-4	94
3	Undecane	156	C11H24	001120-21-4	90
4	Tetradecane	198	C14H30	000629-59-4	90
5	Decane	142	C10H22	000124-18-5	90





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Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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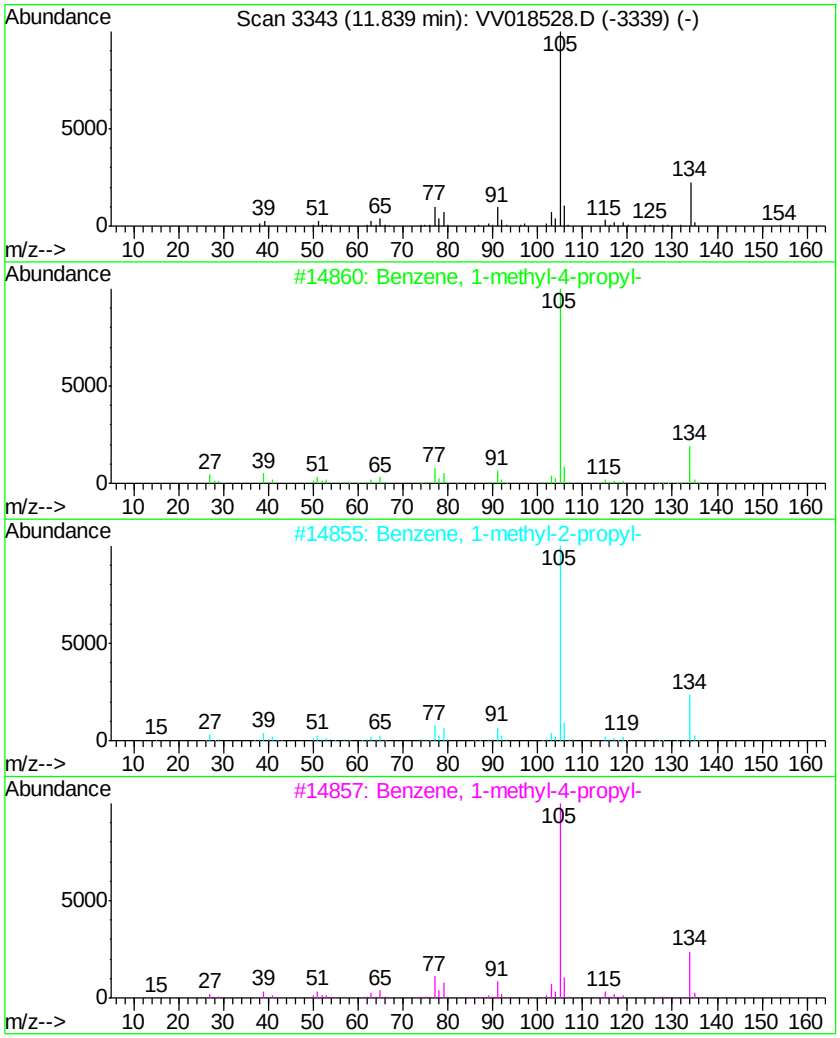
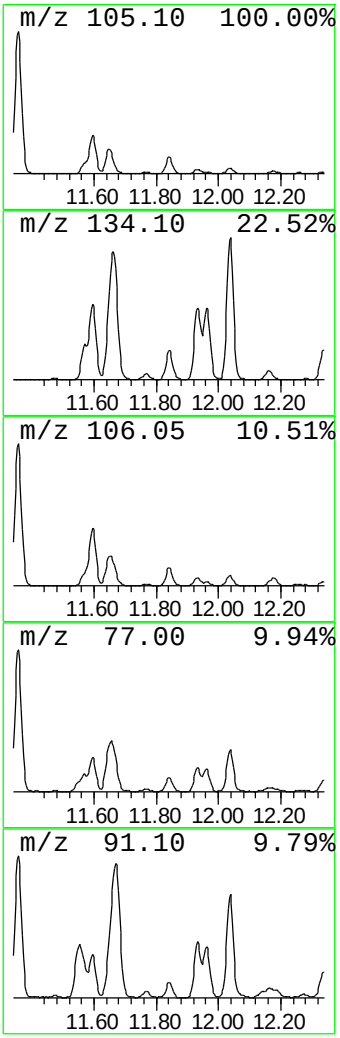
Peak Number 16 Benzene, 1-methyl-4-propyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.84	5.05 ug/L	138253	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
2	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91
3	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	91
5	Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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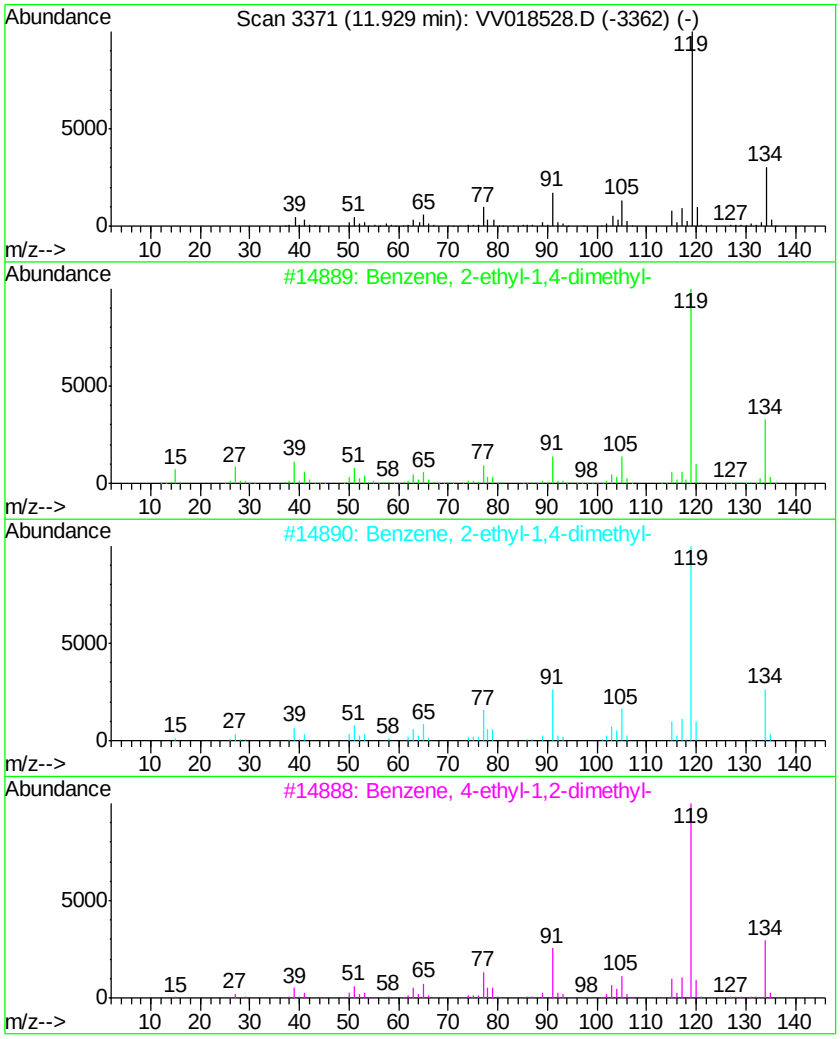
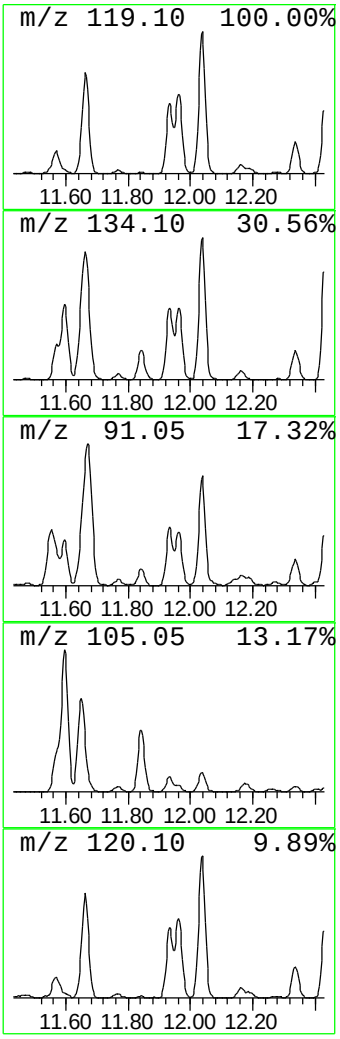
Peak Number 17 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	10.47 ug/L	286691	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	p-Cymene	134	C10H14	000099-87-6	95





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Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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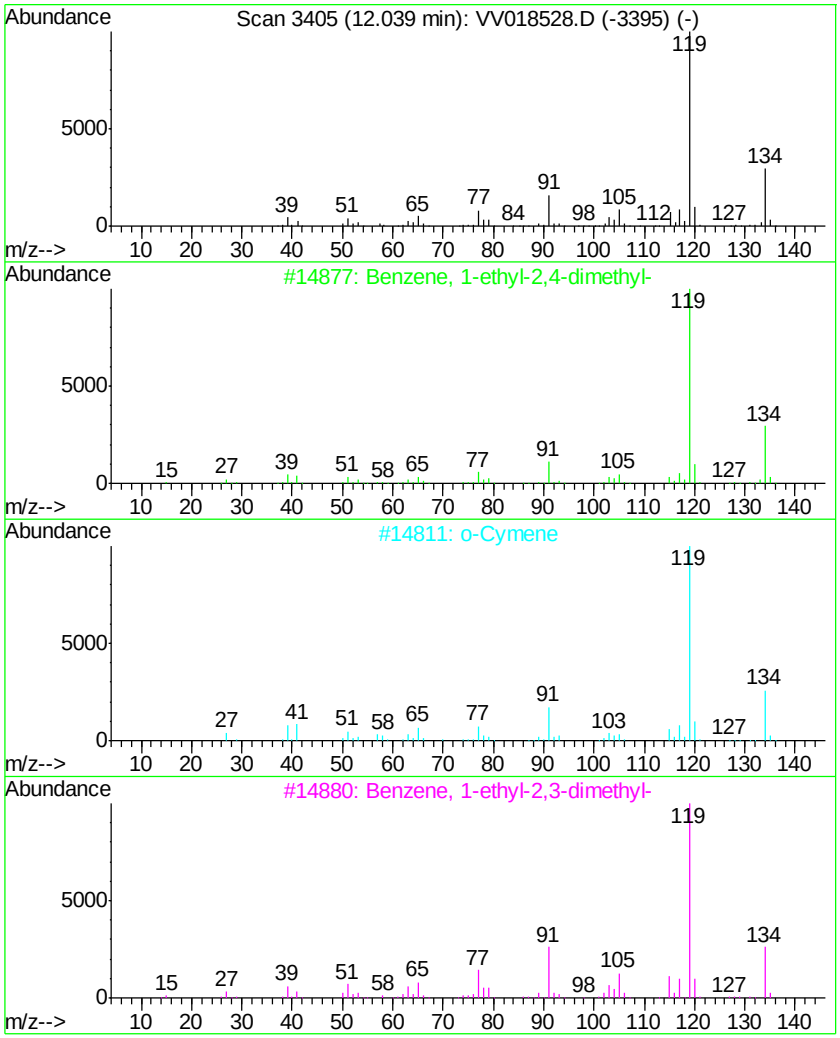
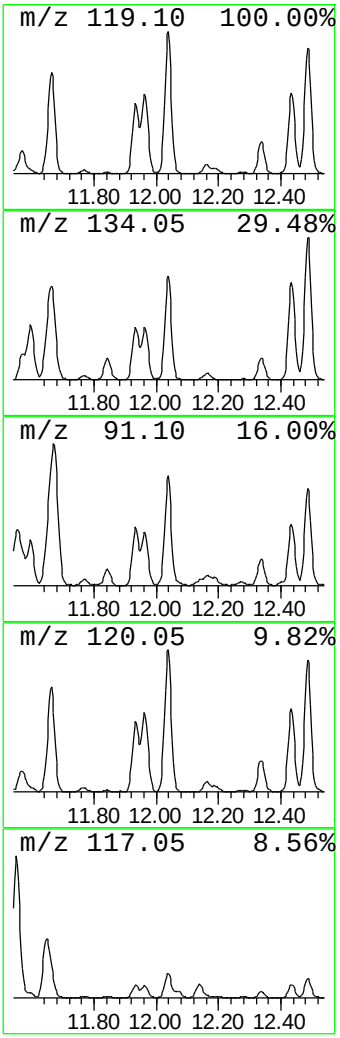
Peak Number 18 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	21.51 ug/L	588769	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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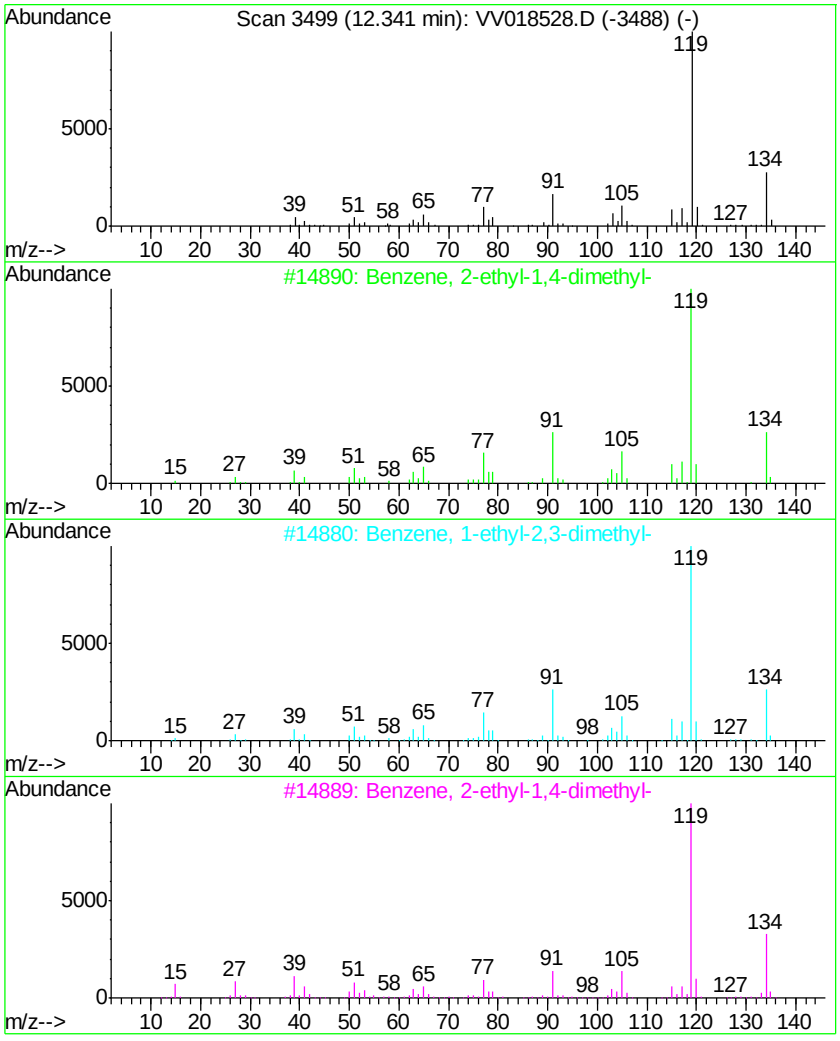
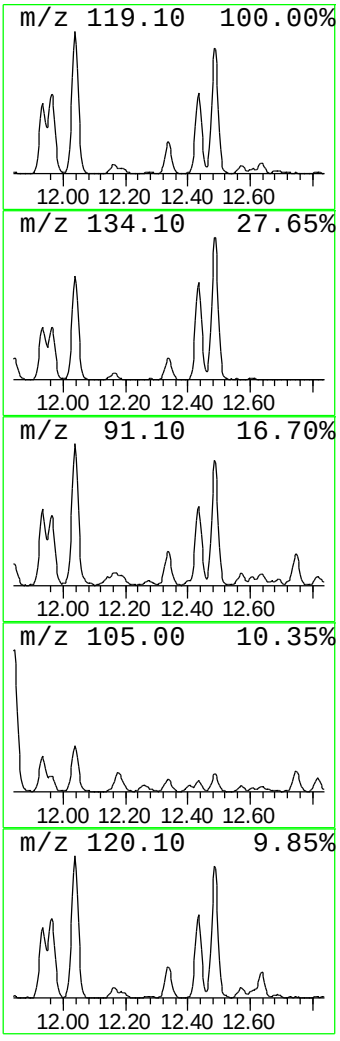
Peak Number 19 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	4.77 ug/L	130513	1,4-Dichlorobenzene-d4	11.27

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





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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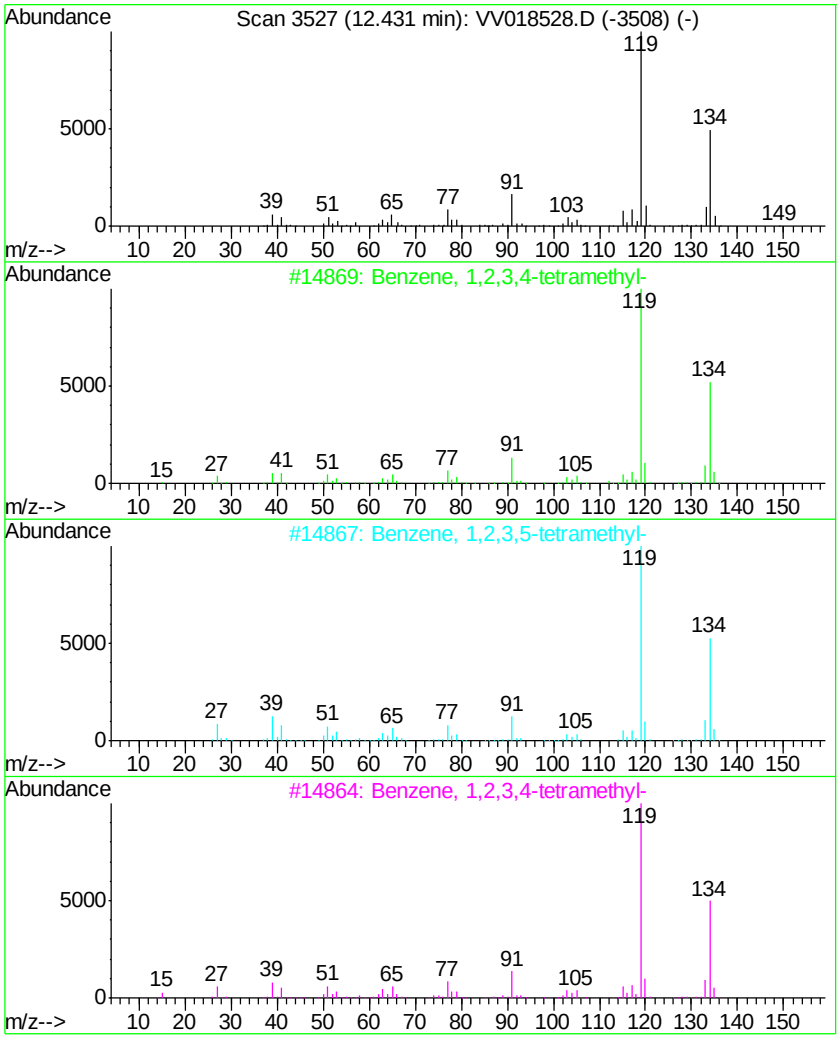
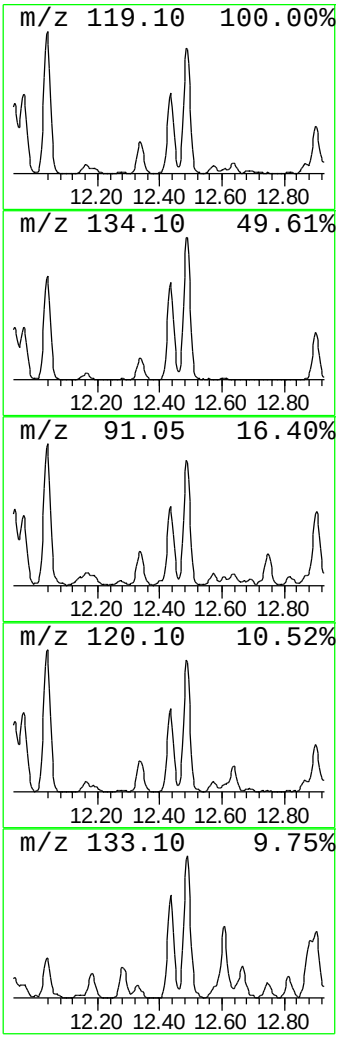
Peak Number 20 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	13.93 ug/L	381361	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	97
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	96





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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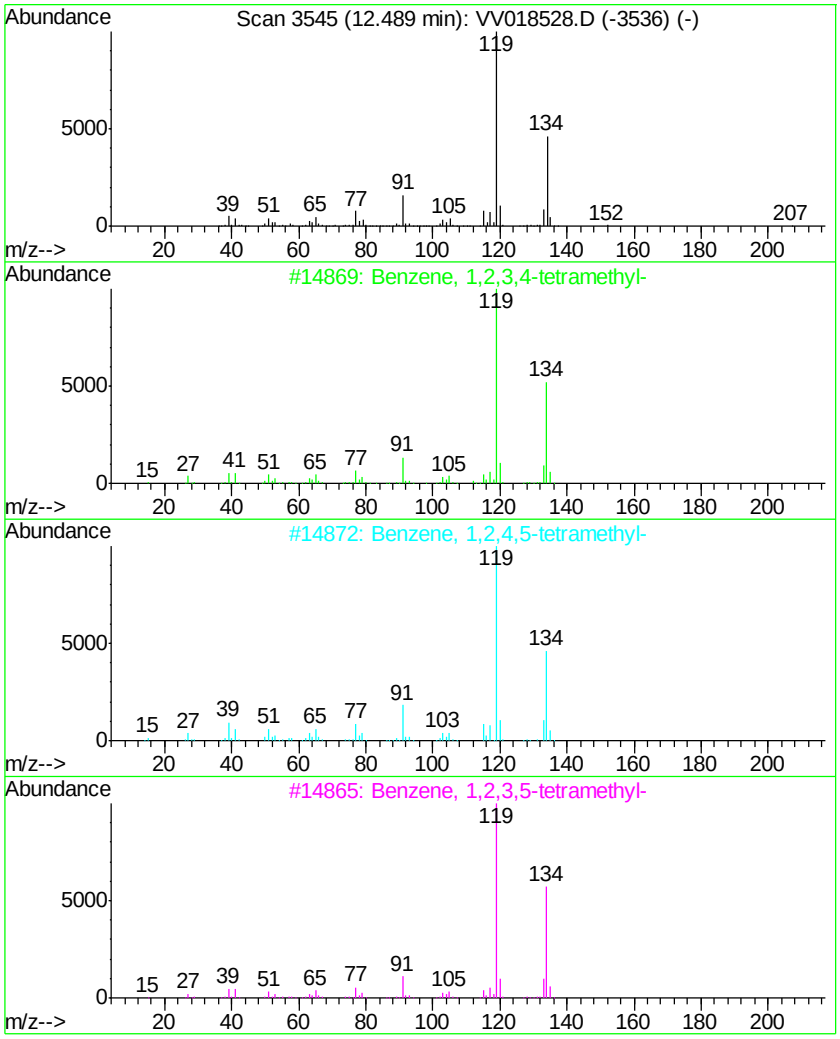
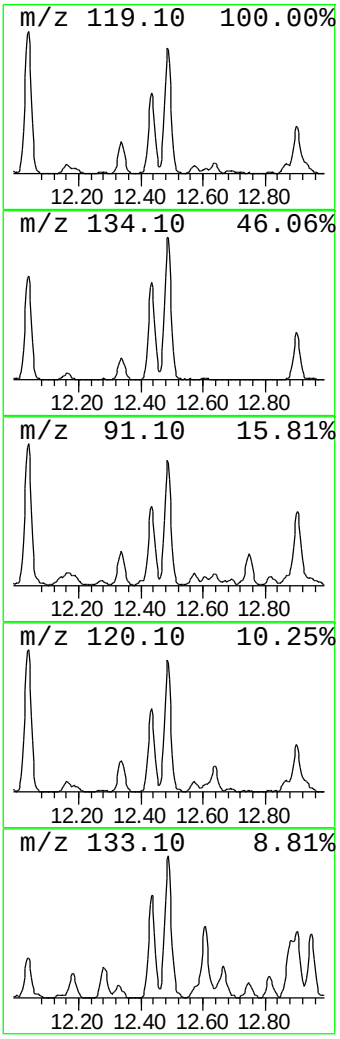
Peak Number 21 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	20.38 ug/L	557984	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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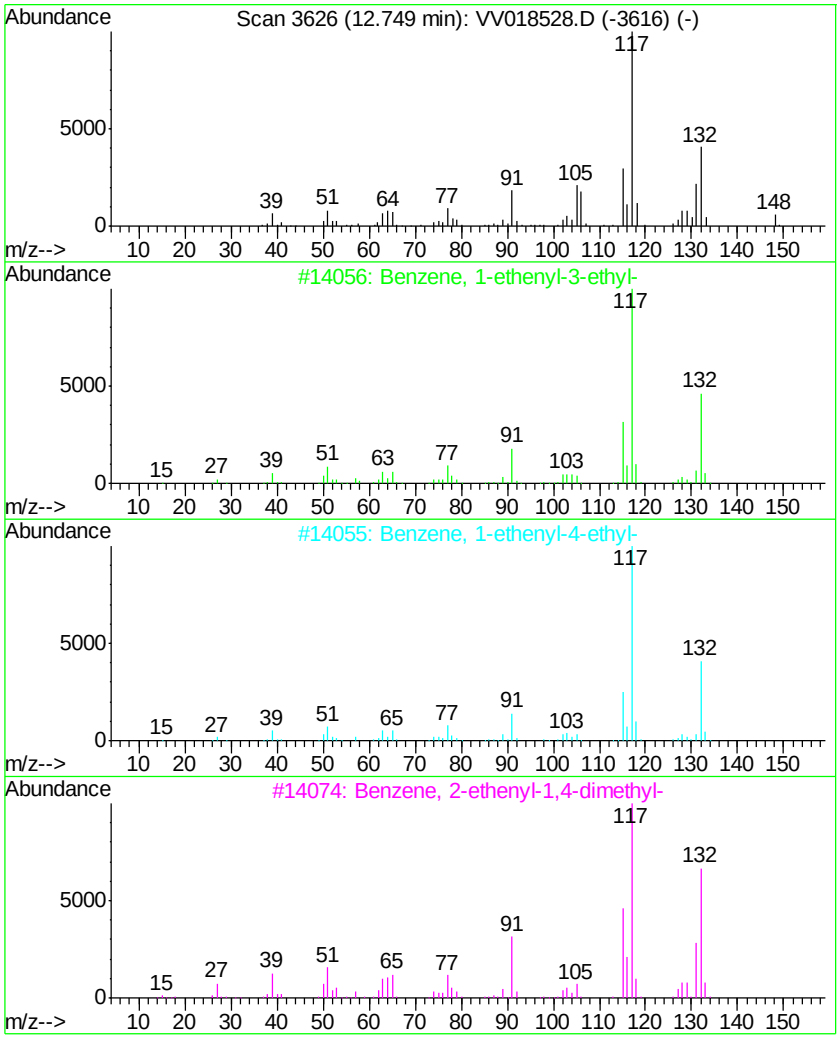
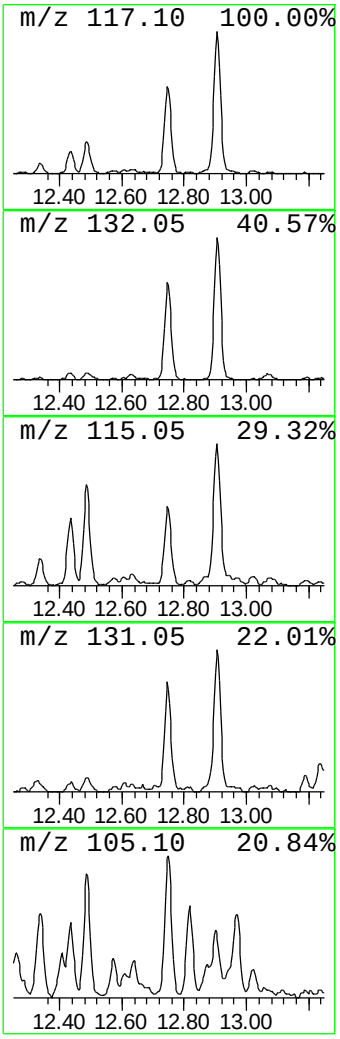
Peak Number 22 Benzene, 1-ethenyl-3-ethyl- Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.75	6.36 ug/L	174153	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	93
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	93
3	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	90
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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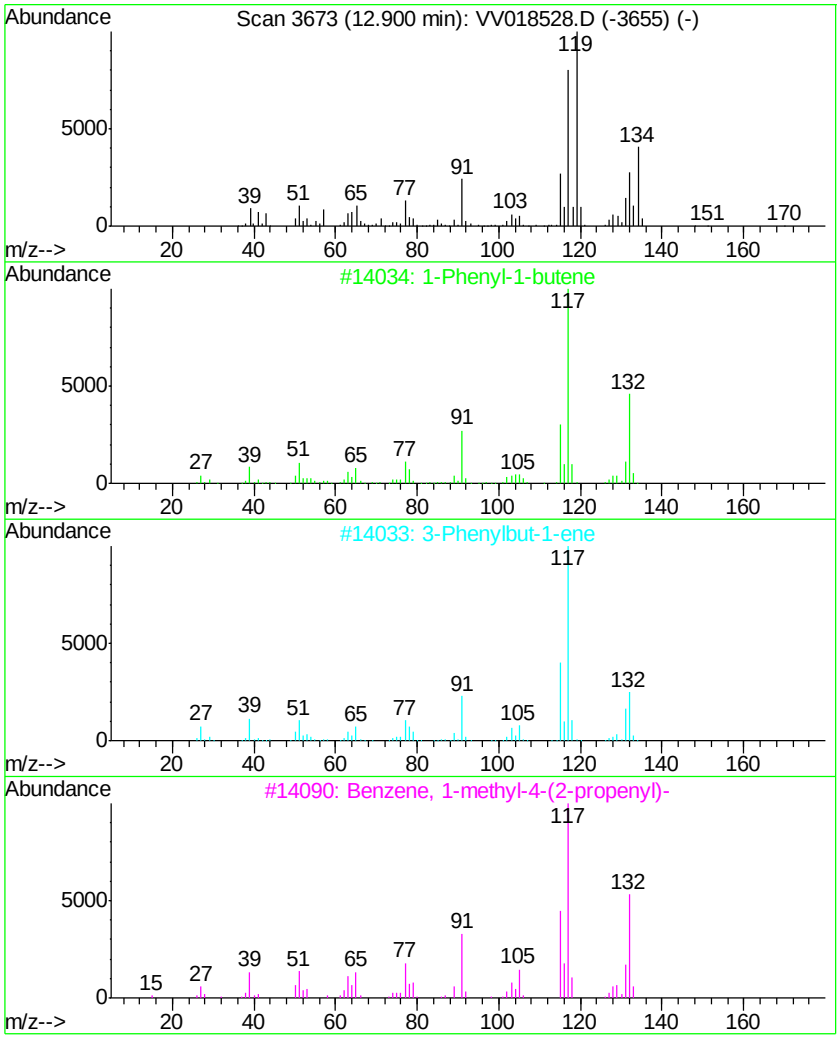
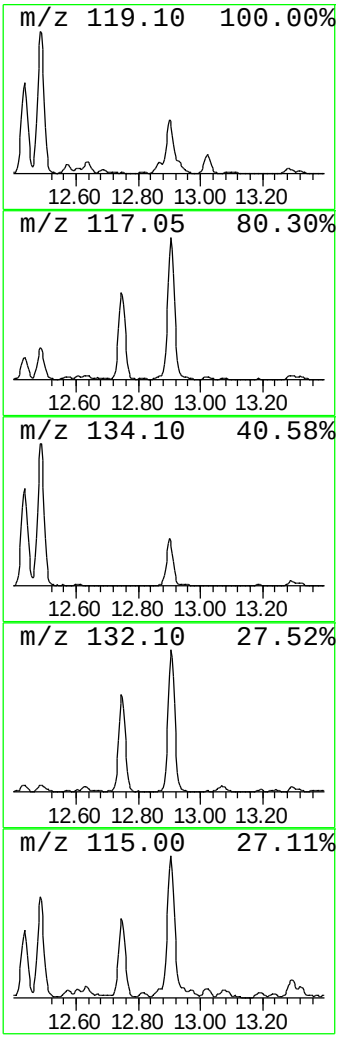
Peak Number 23 1-Phenyl-1-butene Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.90	19.97 ug/L	546760	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	86
3	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
4	p-Cymene	134	C10H14	000099-87-6	50
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	50





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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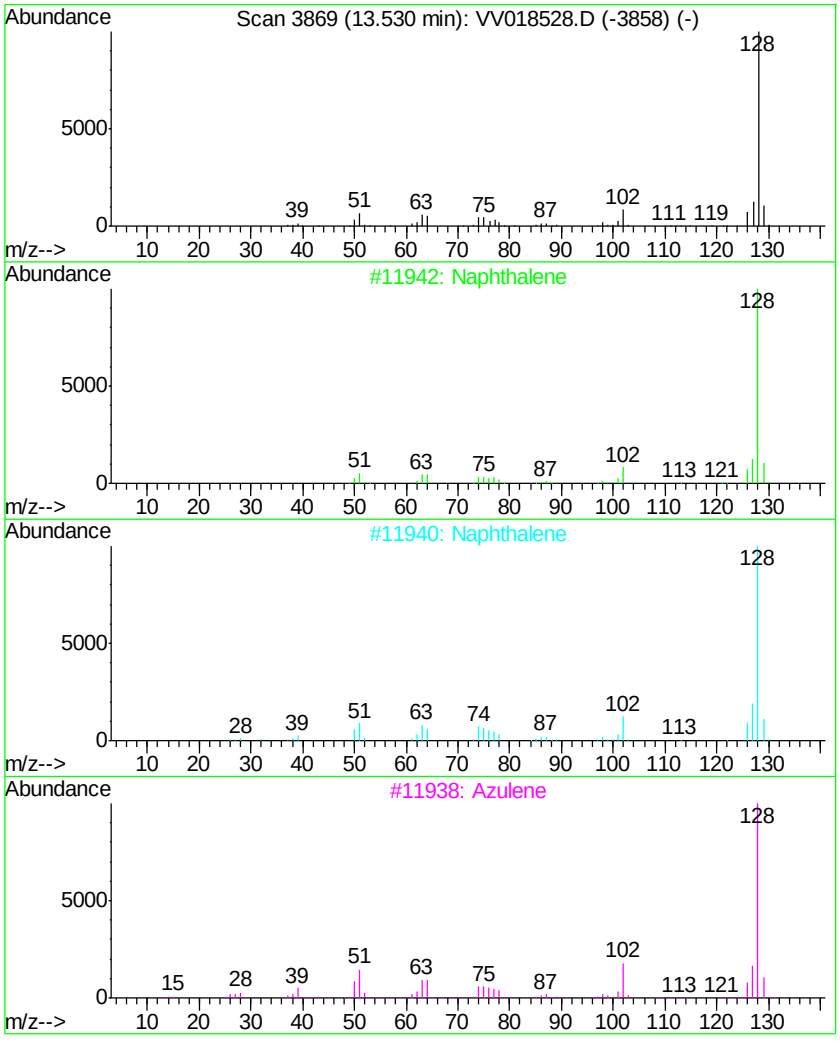
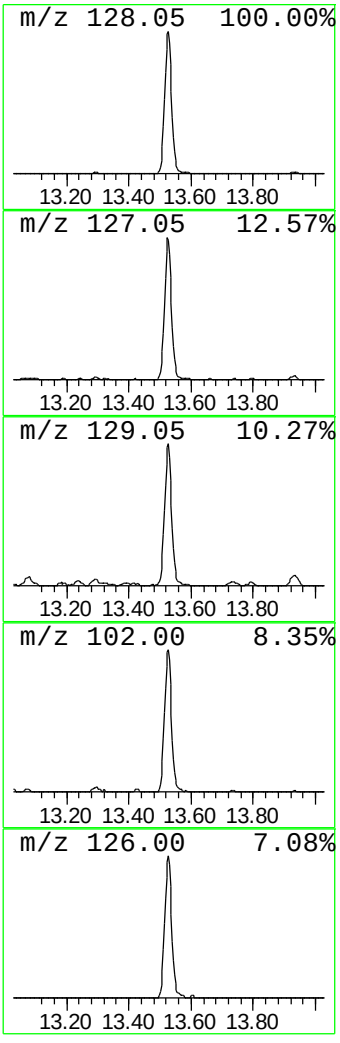
Peak Number 24 Azulene Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.53	12.18 ug/L	333282	1,4-Dichlorobenzene-d4	11.27

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	95
2	Naphthalene	128	C10H8	000091-20-3	94
3	Azulene	128	C10H8	000275-51-4	91
4	Azulene	128	C10H8	000275-51-4	91
5	Azulene	128	C10H8	000275-51-4	90





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018528.D
Acq On : 29 Sep 2020 19:51
Operator : SY/MD
Sample : L4171-14 20X
Misc : 6.13g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 23 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Y4

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: Str...	2.55	42.0	ug/L	859363	1	5.64	1022590	50.0
(DEL) Alkane: Str...	2.78	67.7	ug/L	1384880	1	5.64	1022590	50.0
(DEL) Alkane: Str...	3.06	171.5	ug/L	3506740	1	5.64	1022590	50.0
(DEL) Alkane: Str...	3.57	4.5	ug/L	91993	1	5.64	1022590	50.0
(DEL) Alkane: Cyc...	3.79	48.8	ug/L	998418	1	5.64	1022590	50.0
Benzene, propyl-	10.37	46.3	ug/L	1267440	3	11.27	1368640	50.0
Benzene, 1-ethyl-...	10.47	230.0	ug/L	6296830	3	11.27	1368640	50.0
Mesitylene	10.56	82.4	ug/L	2254990	3	11.27	1368640	50.0
Benzene, 1-ethyl-...	10.76	56.9	ug/L	1558240	3	11.27	1368640	50.0
Benzene, 1,2,3-tr...	10.94	240.5	ug/L	6583270	3	11.27	1368640	50.0
o-Cymene	11.22	3.1	ug/L	84359	3	11.27	1368640	50.0
unknown-01	11.36	51.6	ug/L	1411890	3	11.27	1368640	50.0
Indane	11.55	17.8	ug/L	485808	3	11.27	1368640	50.0
(DEL) Alkane: Str...	11.81	2.9	ug/L	79826	3	11.27	1368640	50.0
Benzene, 1-methyl...	11.84	5.0	ug/L	138253	3	11.27	1368640	50.0
Benzene, 2-ethyl-...	11.93	10.5	ug/L	286691	3	11.27	1368640	50.0
Benzene, 1-ethyl-...	12.04	21.5	ug/L	588769	3	11.27	1368640	50.0
Benzene, 1-ethyl-...	12.34	4.8	ug/L	130513	3	11.27	1368640	50.0
Benzene, 1,2,3,4-...	12.43	13.9	ug/L	381361	3	11.27	1368640	50.0
Benzene, 1,2,4,5-...	12.49	20.4	ug/L	557984	3	11.27	1368640	50.0
Benzene, 1-etheny...	12.75	6.4	ug/L	174153	3	11.27	1368640	50.0
1-Phenyl-1-butene	12.90	20.0	ug/L	546760	3	11.27	1368640	50.0
Azulene	13.53	12.2	ug/L	333282	3	11.27	1368640	50.0