LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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	1.312	62	69	87	rVB	137887	165278	0.89%	0.345%
2	1.576	143	151	168	rBV	92266	139689	0.75%	0.292%
3	2.100	303	314	327	rBV	272369	393455	2.11%	0.822%
4	2.528	420	447	470	rBV	1894779	3737735	20.09%	7.804%
5	2.627	471	478	495	rVB	8421	18805	0.10%	0.039%
6	2.762	502	520	546	rBV	2751393	5506779	29.60%	11.498%
7	3.045	588	608	638	rBV	9373526	18605587	100.00%	38.849%
8	3.550	743	765	785	rBV	556652	1454425	7.82%	3.037%
9	3.679	787	805	820	rVV	476926	1242894	6.68%	2.595%
10	3.772	820	834	854	rVV	473462	1190736	6.40%	2.486%
11	3.891	857	871	874	rVV2	26902	51105	0.27%	0.107%
12	3.930	875	883	929	rVB2	54280	197059	1.06%	0.411%
13	4.370	1003	1020	1052	rBV	206813	557599	3.00%	1.164%
14	4.692	1103	1120	1135	rBV3	45170	115307	0.62%	0.241%
15	4.785	1136	1149	1167	rVB3	19837	51998	0.28%	0.109%
16	4.926	1174	1193	1220	rBV3	20581	54835	0.29%	0.114%
17	5.068	1220	1237	1264	rBV2	539622	1330226	7.15%	2.778%
18	5.508	1360	1374	1394	rBV2	21171	44720	0.24%	0.093%
19	5.637	1401	1414	1451	rBV	434672	906572	4.87%	1.893%
20	5.897	1479	1495	1496	rBV2	17354	24950	0.13%	0.052%
21	6.093	1542	1556	1584	rBV	280490	630551	3.39%	1.317%
22	6.216	1585	1594	1604	rVV6	10484	20553	0.11%	0.043%
23	6.653	1714	1730	1752	rBV6	10498	28880	0.16%	0.060%
24	6.859	1784	1794	1803	rBV4	8897	17988	0.10%	0.038%
25	6.936	1808	1818	1826	rBV2	37368	61942	0.33%	0.129%
26	7.010	1828	1841	1861	rVV	161175	334716	1.80%	0.699%
27	7.097	1861	1868	1887	rVB2	34691	69260	0.37%	0.145%
28	7.187	1887	1896	1913	rVB9	7904	20919	0.11%	0.044%
29	7.283	1913	1926	1932	rBV2	76688	143522	0.77%	0.300%
30	7.338	1932	1943	1968	rVB	639553	1178397	6.33%	2.461%
31	7.463	1971	1982	1993	rBV3	25894	50560	0.27%	0.106%
32	7.585	2011	2020	2030	rBV	82271	130791	0.70%	0.273%
33	7.646	2030	2039	2043	rBV	100413	164909	0.89%	0.344%
34	7.807	2074	2089	2100	rBV2	33321	61745	0.33%	0.129%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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	7.997	2135	2148	2160	rVB3	14108	25645	0.14%	0.054%
36	8.142	2169	2193	2212	rBV	162861	583932	3.14%	1.219%
37	8.309	2236	2245	2255	rVB2	35426	58168	0.31%	0.121%
38	8.370	2255	2264	2275	rBV2	26554	45277	0.24%	0.095%
39	8.685	2354	2362	2372	rVB3	15132	25950	0.14%	0.054%
40	8.807	2390	2400	2409	rBV3	10773	18702	0.10%	0.039%
41	8.875	2409	2421	2441	rBV	617283	1157819	6.22%	2.418%
42	9.225	2522	2530	2549	rVB4	18940	34892	0.19%	0.073%
43	9.495	2600	2614	2632	rVB10	9927	19773	0.11%	0.041%
44	9.730	2680	2687	2697	rVB4	14012	21295	0.11%	0.044%
45	9.817	2698	2714	2724	rBV8	10674	23245	0.12%	0.049%
46	10.106	2791	2804	2809	rVV3	13593	24982	0.13%	0.052%
47	10.145	2809	2816	2826	rVB5	22068	38952	0.21%	0.081%
48	10.248	2826	2848	2864	rBV	359086	638634	3.43%	1.333%
49	10.476	2909	2919	2930	rBV2	25624	54484	0.29%	0.114%
50	10.598	2942	2957	2979	rVB4	73494	149396	0.80%	0.312%
51	10.900	3043	3051	3057	rBV3	34207	51807	0.28%	0.108%
52	10.939	3057	3063	3073	rVB	38191	57544	0.31%	0.120%
53	11.058	3089	3100	3103	rBV4	12022	19511	0.10%	0.041%
54	11.084	3105	3108	3112	rVV3	15333	18745	0.10%	0.039%
55	11.116	3114	3118	3127	rVV5	23198	37690	0.20%	0.079%
56	11.174	3128	3136	3143	rVV5	29387	54883	0.29%	0.115%
57	11.215	3143	3149	3157	rVV7	28496	50769	0.27%	0.106%
58	11.273	3157	3167	3187	rVV	806904	1358571	7.30%	2.837%
59	11.357	3187	3193	3199	rVV4	27272	46411	0.25%	0.097%
60	11.396	3199	3205	3215	rVV2	29985	56770	0.31%	0.119%
61	11.441	3215	3219	3226	rVV7	12916	20842	0.11%	0.044%
62	11.486	3226	3233	3243	rVB3	25059	37875	0.20%	0.079%
63	11.572	3251	3260	3263	rBV	24006	35443	0.19%	0.074%
64	11.598	3264	3268	3274	rVV	39224	54153	0.29%	0.113%
65	11.650	3274	3284	3301	rVB	783567	1290911	6.94%	2.695%
66	11.810	3325	3334	3341	rBV2	91083	129799	0.70%	0.271%
67	11.939	3365	3374	3377	rBV2	33408	50819	0.27%	0.106%
68	11.965	3377	3382	3390	rVV2	47813	71588	0.38%	0.149%
69	12.039	3391	3405	3419	rVB	55167	107146	0.58%	0.224%
70	12.145	3430	3438	3441	rBV4	14761	20762	0.11%	0.043%
71	12.187	3444	3451	3468	rVB5	33178	76125	0.41%	0.159%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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	12.283	3468	3481	3489	rBV6	34399	70262	0.38%	0.147%
73	12.331	3489	3496	3505	rVB7	14971	29426	0.16%	0.061%
74	12.392	3505	3515	3521	rBV5	32877	62488	0.34%	0.130%
75	12.434	3522	3528	3537	rVB2	45932	71928	0.39%	0.150%
76	12.489	3537	3545	3553	rBV	79905	122171	0.66%	0.255%
77	12.572	3563	3571	3577	rBV2	35155	51792	0.28%	0.108%
78	12.611	3577	3583	3587	rVV2	22354	28214	0.15%	0.059%
79	12.749	3620	3626	3638	rVB2	32161	46960	0.25%	0.098%
80	12.817	3640	3647	3655	rBV2	22976	32314	0.17%	0.067%
81	12.871	3655	3664	3668	rBV3	31592	50077	0.27%	0.105%
82	12.907	3669	3675	3704	rVB6	73486	181213	0.97%	0.378%
83	13.022	3705	3711	3718	rBV	36475	47579	0.26%	0.099%
84	13.061	3718	3723	3729	rBV3	13995	18807	0.10%	0.039%
85	13.186	3754	3762	3770	rBV4	17191	24394	0.13%	0.051%
86	13.286	3782	3793	3812	rBV2	698788	1109742	5.96%	2.317%
87	13.424	3830	3836	3849	rVB	25578	42842	0.23%	0.089%
88	13.502	3850	3860	3874	rVV	20140	55359	0.30%	0.116%
89	13.566	3874	3880	3889	rVB9	15975	24652	0.13%	0.051%
90	13.640	3895	3903	3908	rBV6	15322	26472	0.14%	0.055%
91	13.768	3934	3943	3957	rVB	136115	227058	1.22%	0.474%
92	13.929	3985	3993	4007	rVB5	27183	57017	0.31%	0.119%
93	14.125	4042	4054	4068	rBV4	25618	63722	0.34%	0.133%
94	14.257	4088	4095	4103	rBV2	14827	21051	0.11%	0.044%
95	14.457	4150	4157	4169	rVV6	12932	23725	0.13%	0.050%
96	14.849	4272	4279	4290	rVB10	12146	25559	0.14%	0.053%
97	15.614	4514	4517	4528	rVB6	16461	21633	0.12%	0.045%
98	15.694	4533	4542	4554	rBV3	18593	36195	0.19%	0.076%
99	15.839	4579	4587	4594	rBV2	18482	26749	0.14%	0.056%
100	15.961	4618	4625	4634	rVB2	11487	19334	0.10%	0.040%

Sum of corrected areas: 47892532

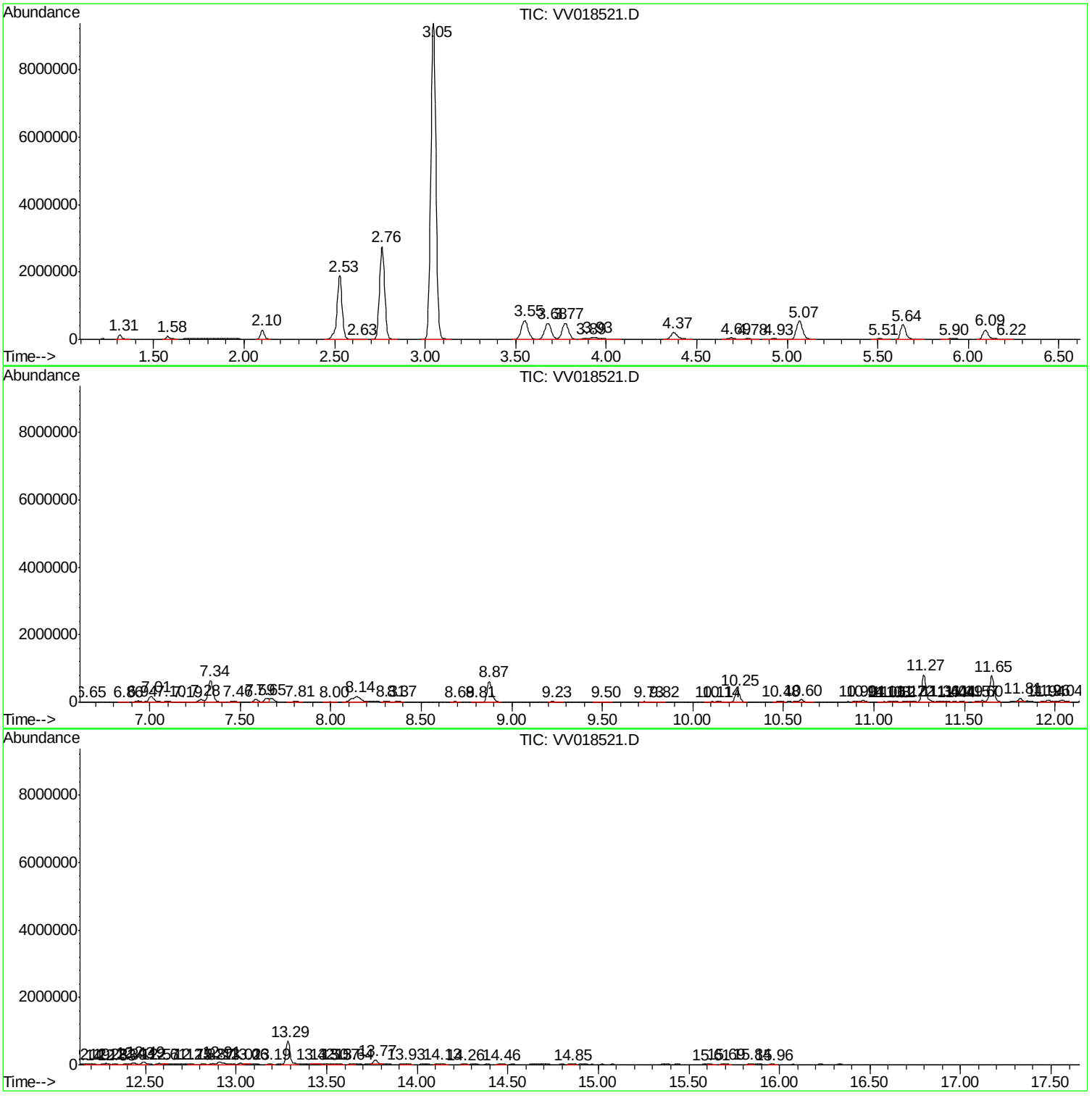
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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 : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

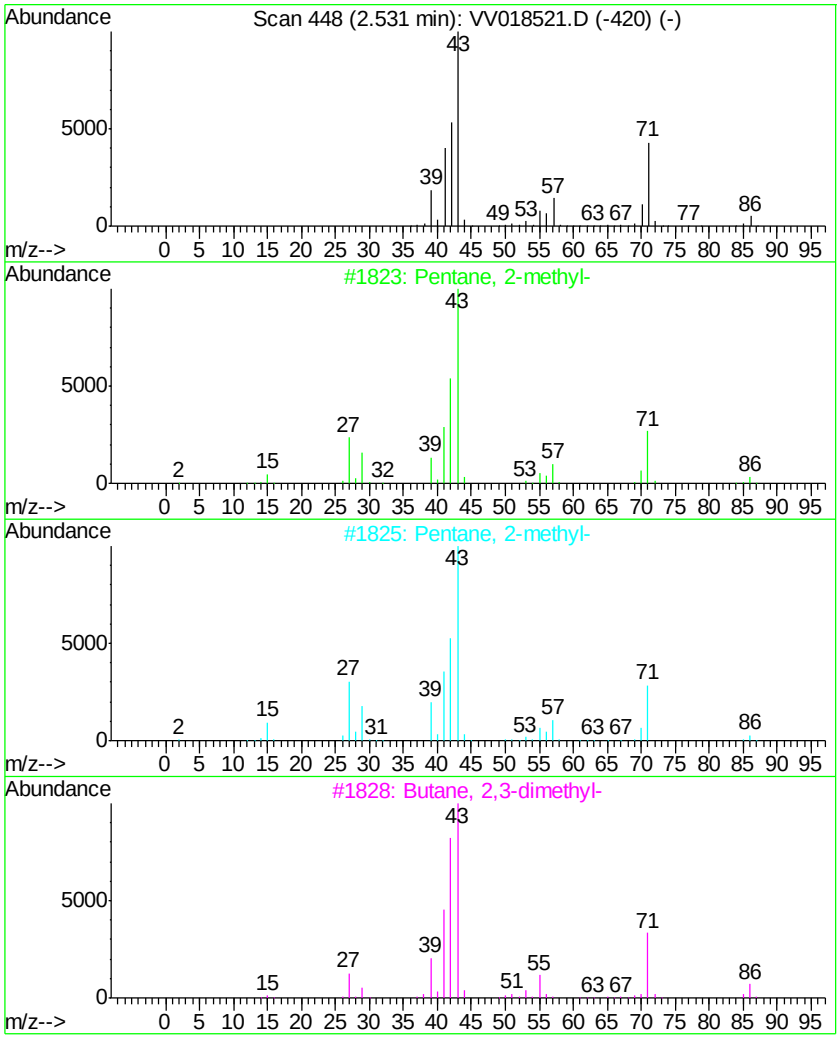
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

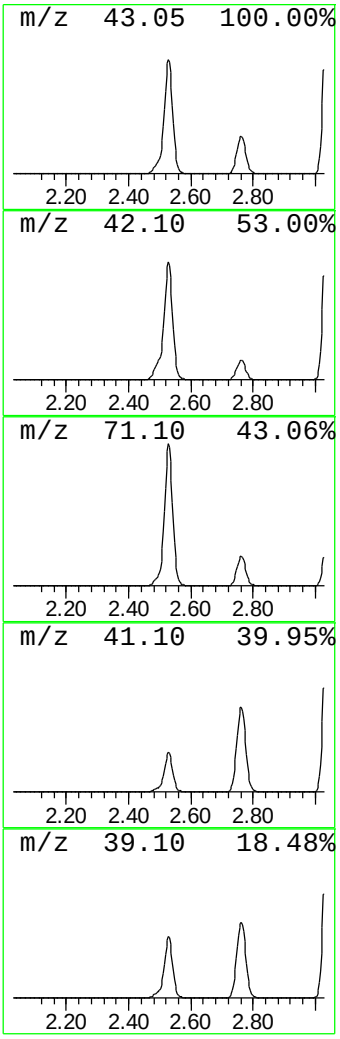
****

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.53	206.15 ug/L	3737740	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 : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

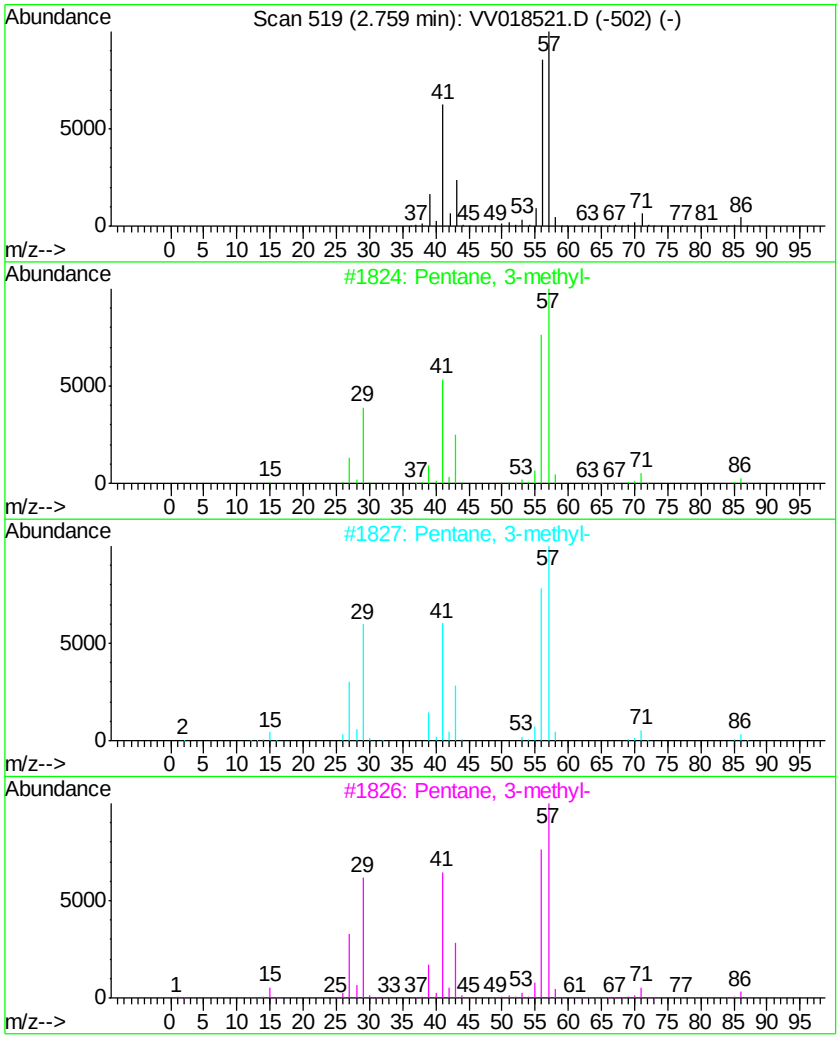
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

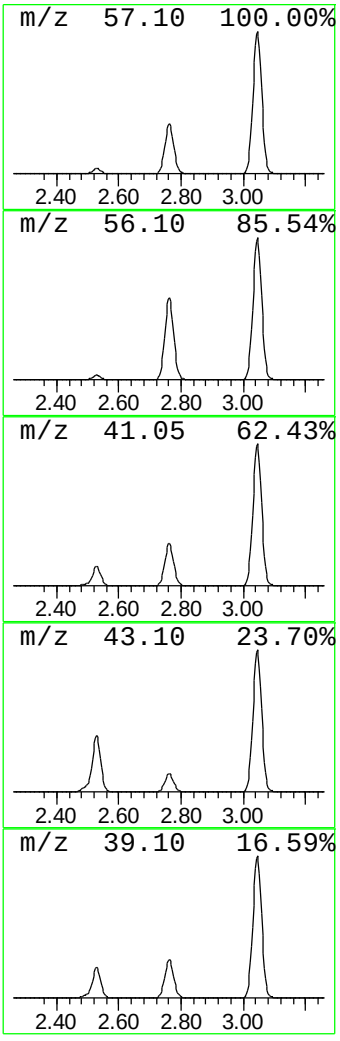
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.76	303.71 ug/L	5506780	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	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	43





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

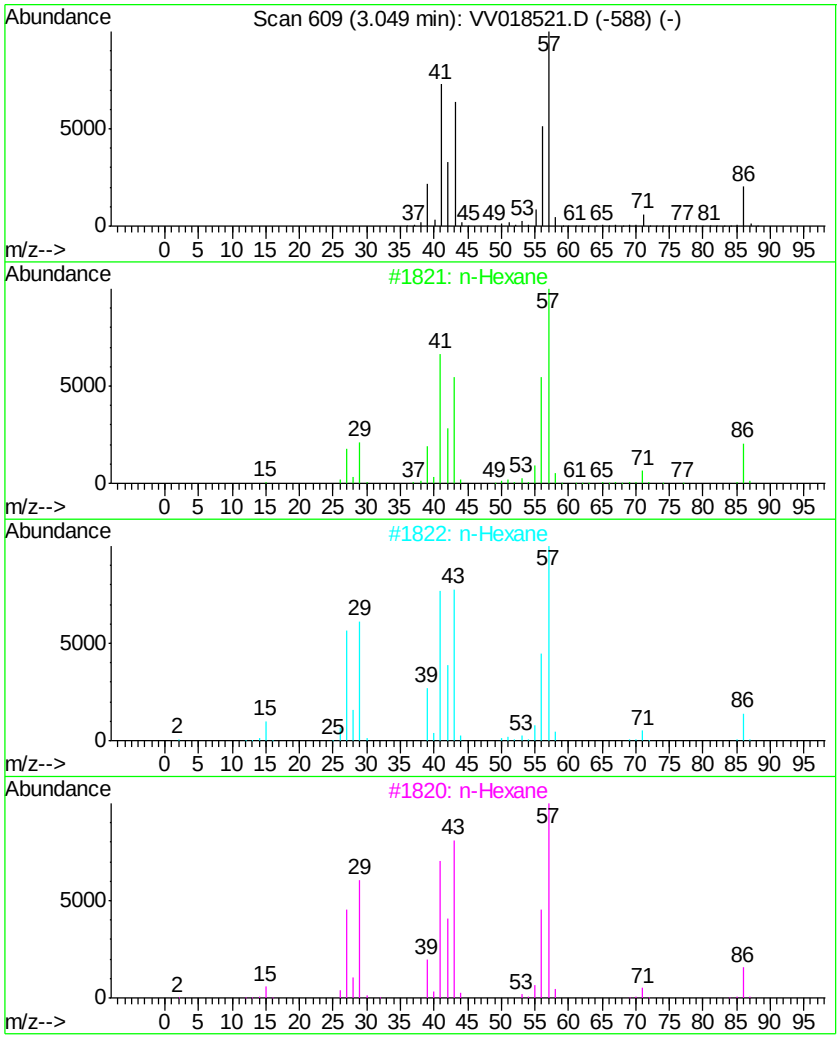
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

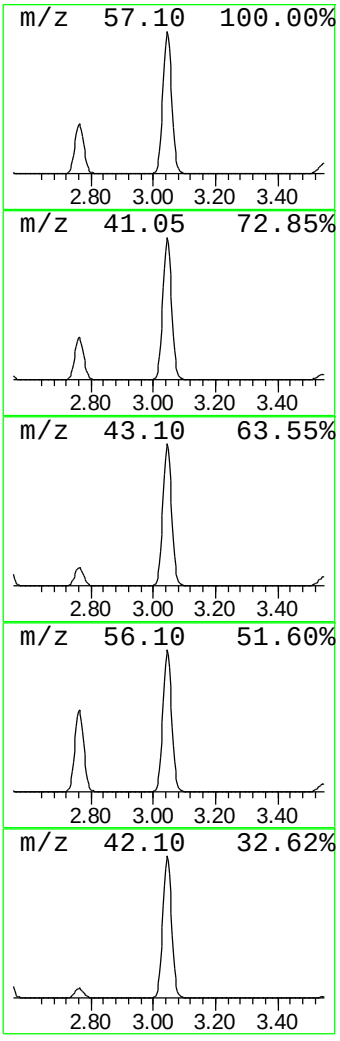
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	1026.15 ug/L	18605600	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	72
3	n-Hexane	86	C6H14	000110-54-3	59
4	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	47
5	Propanal, 2,2-dimethyl-	86	C5H10O	000630-19-3	43





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

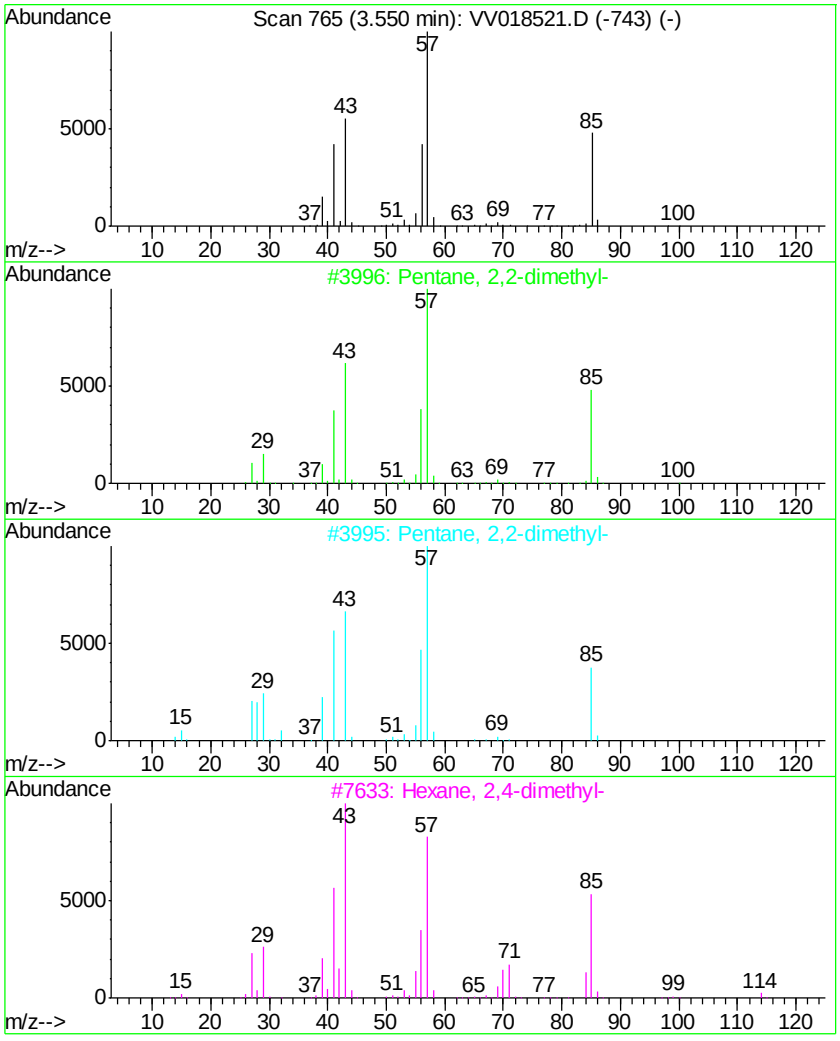
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

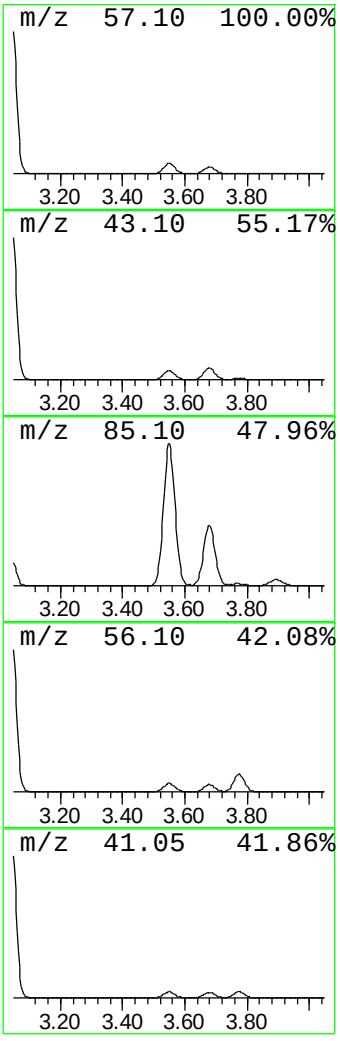
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.55	80.22 ug/L	1454430	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	90
2	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	83
3	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	72
4	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
5	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

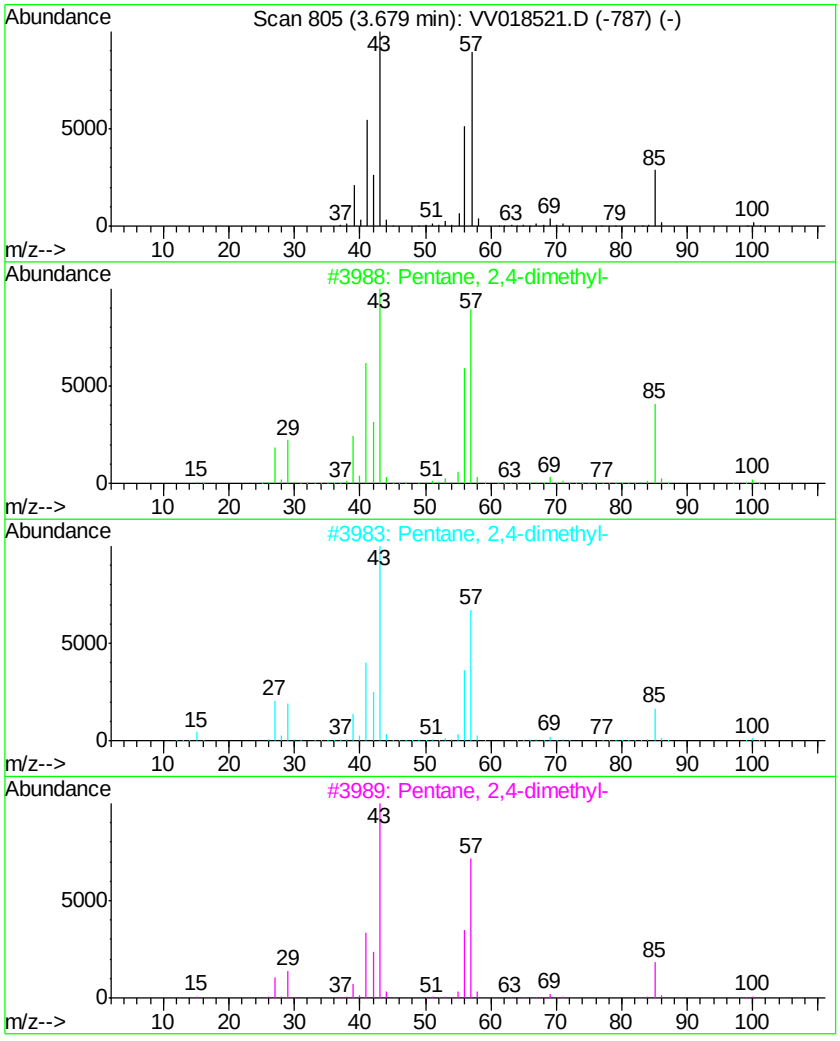
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

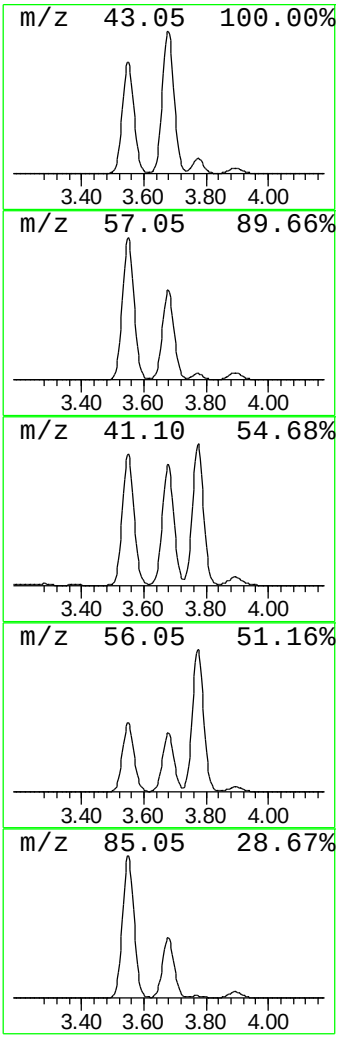
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.68	68.55 ug/L	1242890	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	94
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
3	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

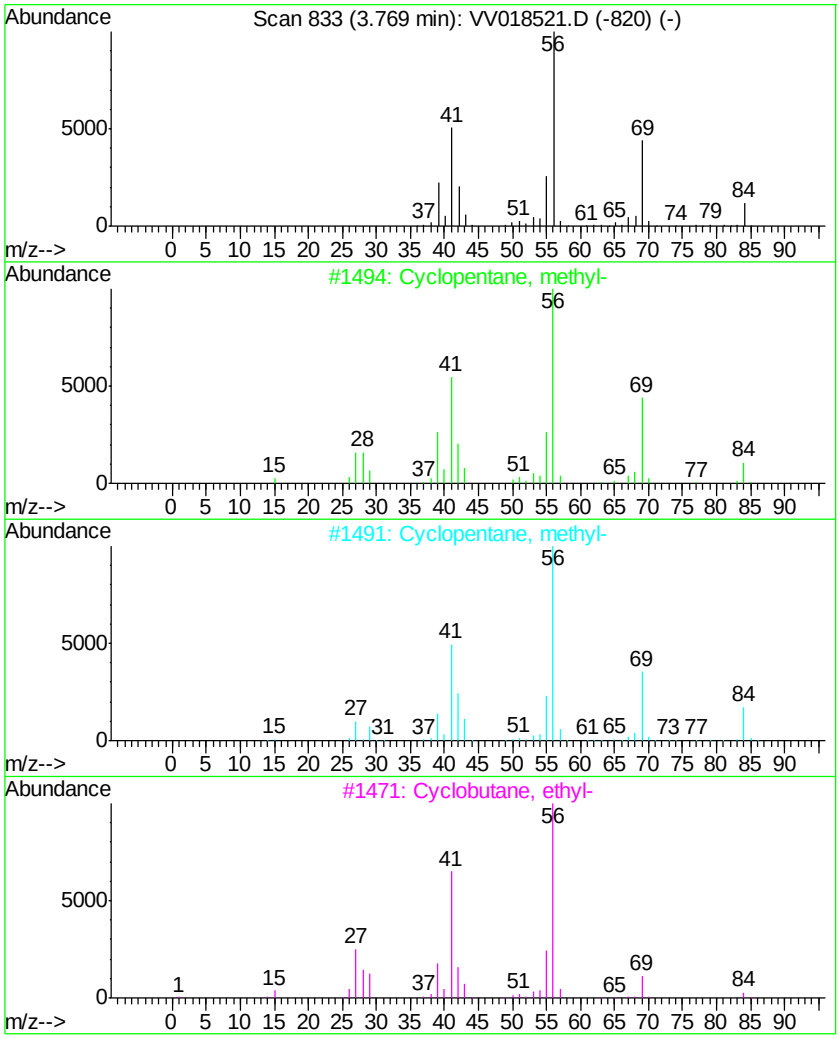
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

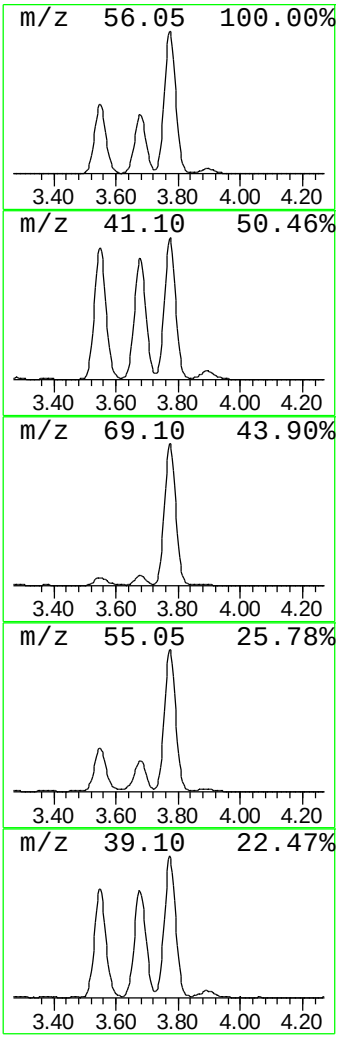
****

Peak Number 6 (DEL) Alkane: Cyclic3.77 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.77	65.67 ug/L	1190740	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	78
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5	Cyclopropane, propyl-	84	C6H12	002415-72-7	56





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

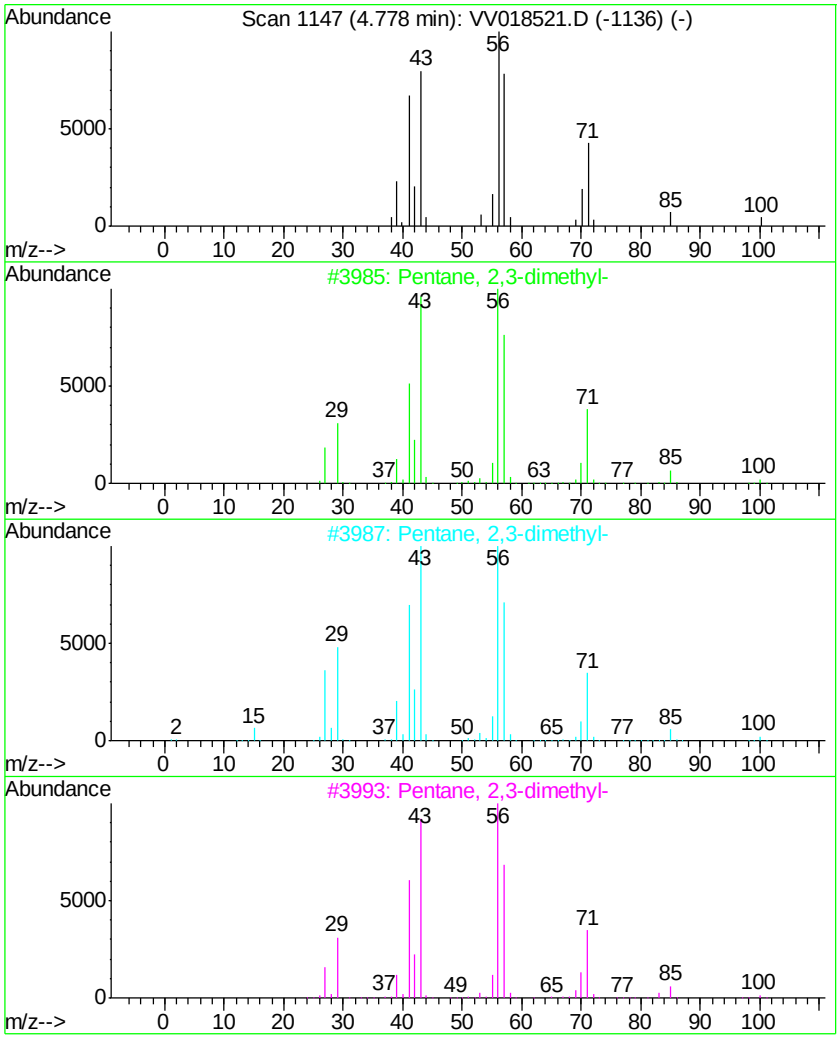
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

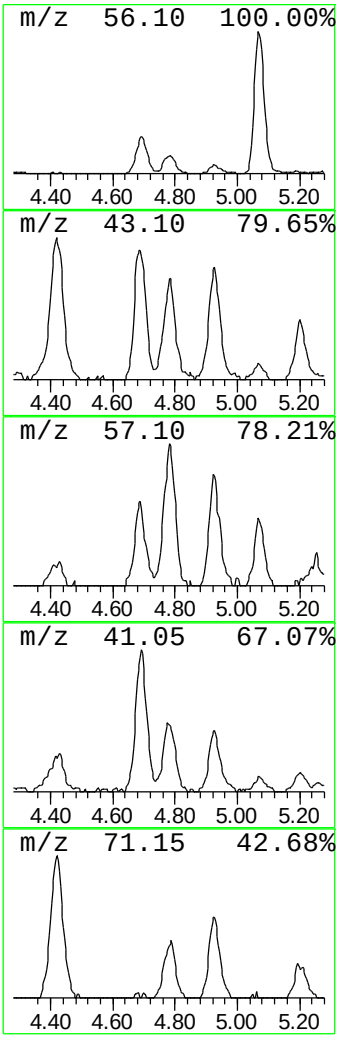
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.78	2.87 ug/L	51998	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	83
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	74
5	Aziridine, 2-methyl-	57	C3H7N	000075-55-8	43





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

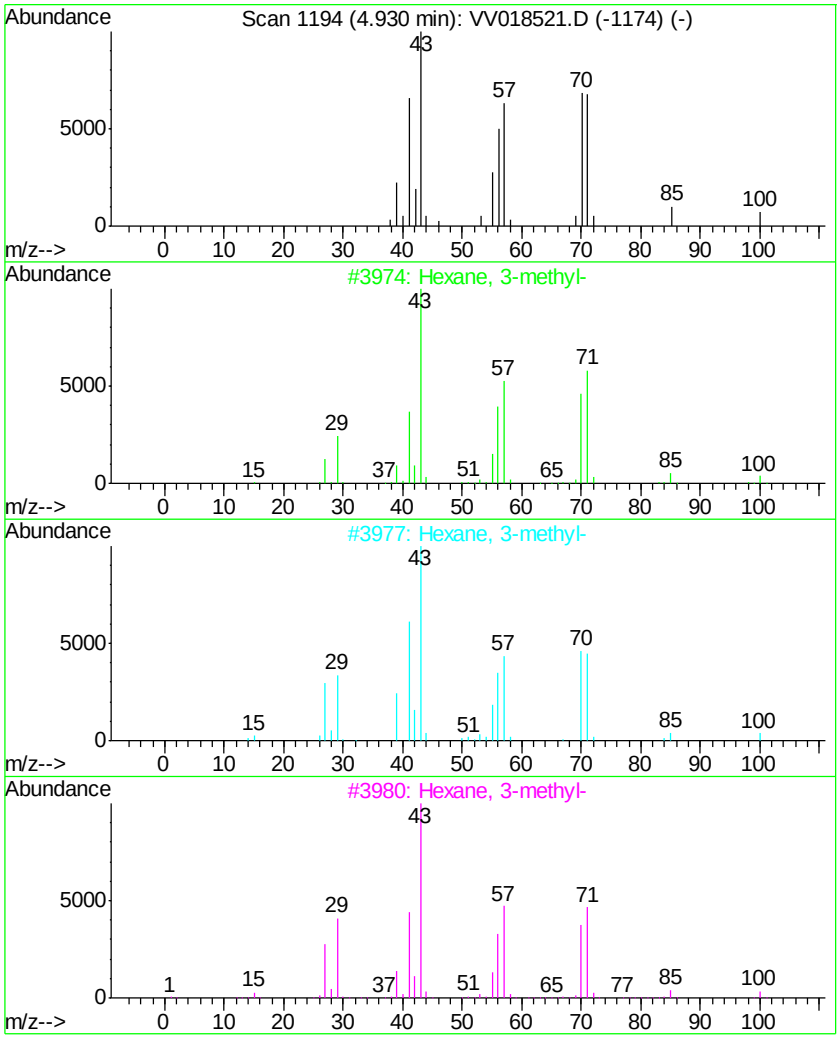
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

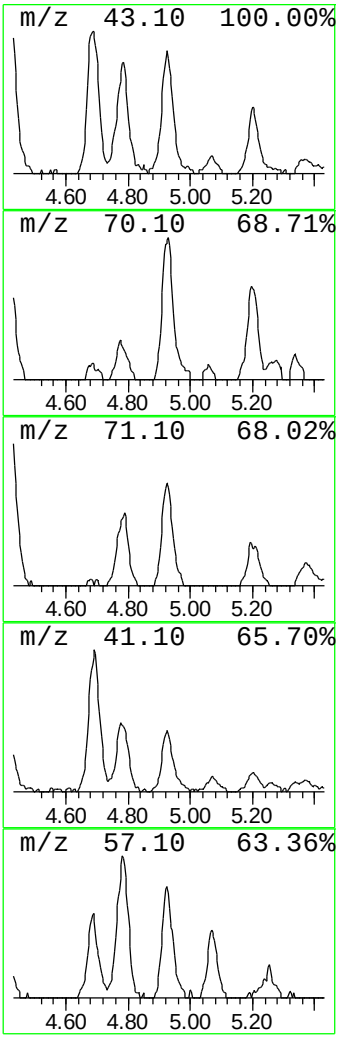
****

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.93	3.02 ug/L	54835	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
2	Hexane, 3-methyl-	100	C7H16	000589-34-4	91
3	Hexane, 3-methyl-	100	C7H16	000589-34-4	87
4	Heptane	100	C7H16	000142-82-5	80
5	Heptane	100	C7H16	000142-82-5	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

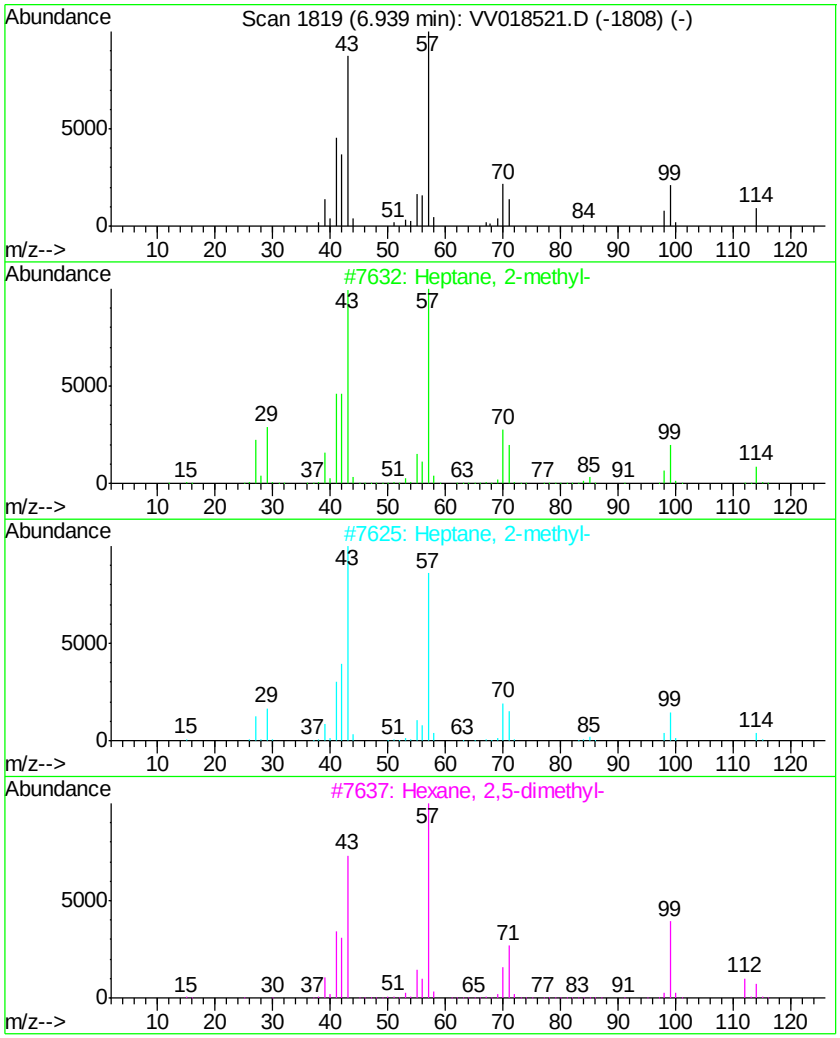
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

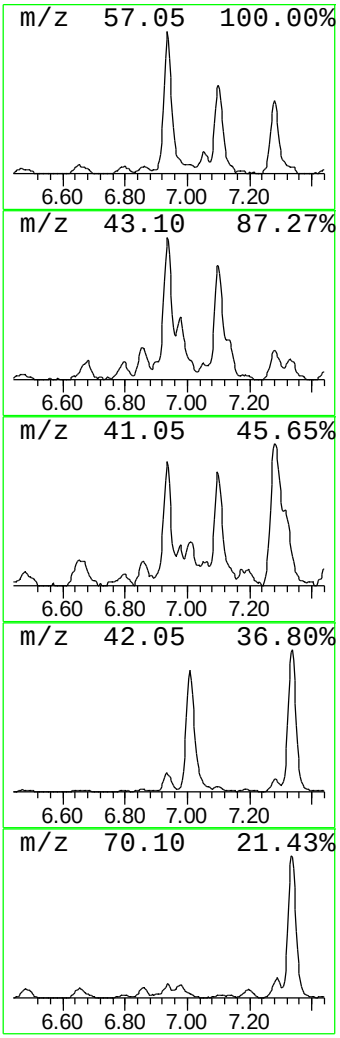
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.94	3.42 ug/L	61942	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 2-methyl-	114	C8H18	000592-27-8	95
2	Heptane, 2-methyl-	114	C8H18	000592-27-8	78
3	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	72
4	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	64
5	Heptane, 2,3,6-trimethyl-	142	C10H22	004032-93-3	53





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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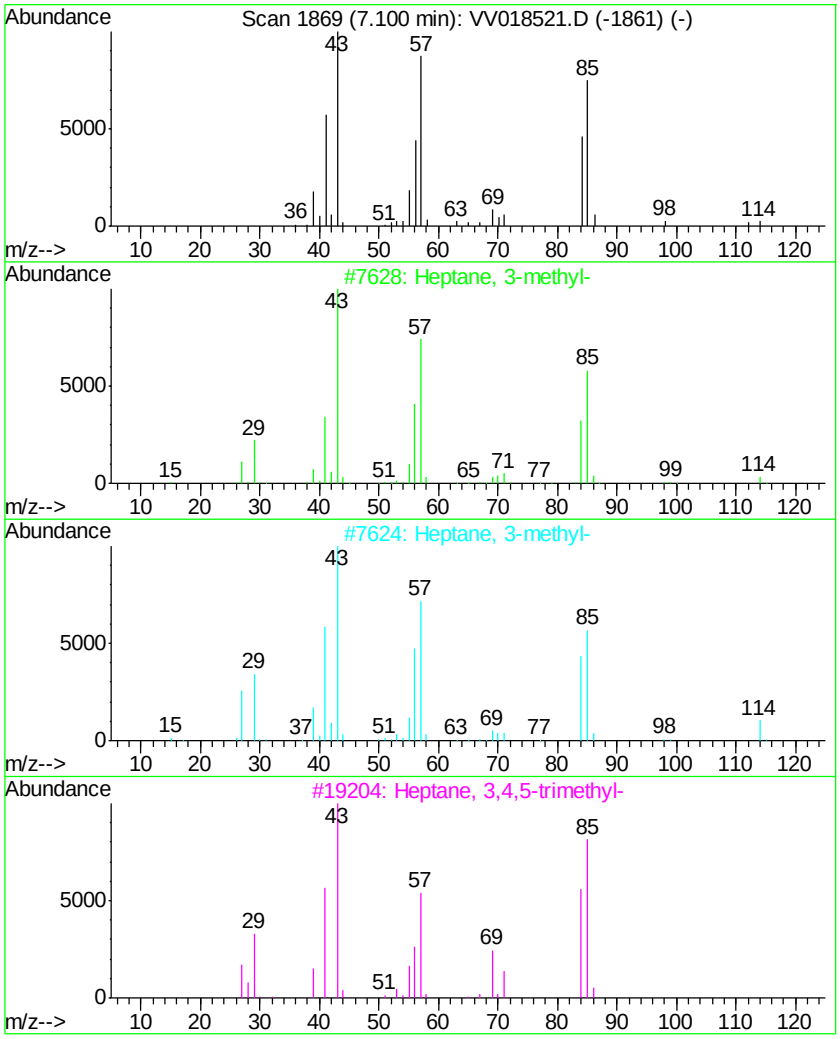
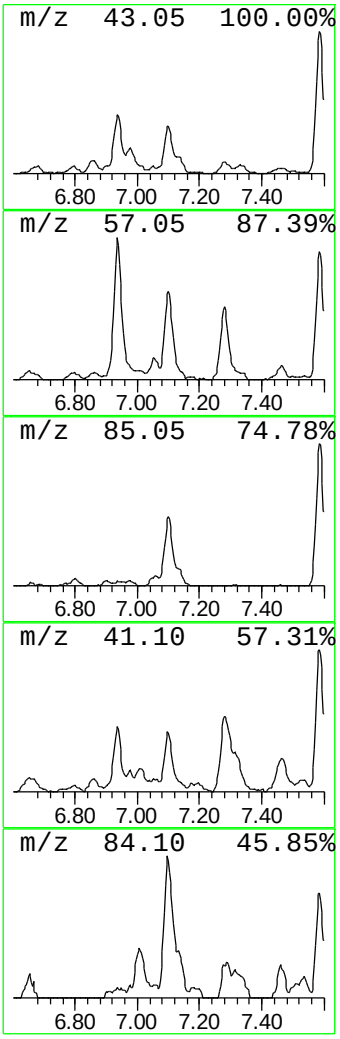
Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.10	3.82 ug/L	69260	1,4-Difluorobenzene	5.64

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2	Heptane, 3-methyl-	114	C8H18	000589-81-1	86
3	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	64
4	Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	64
5	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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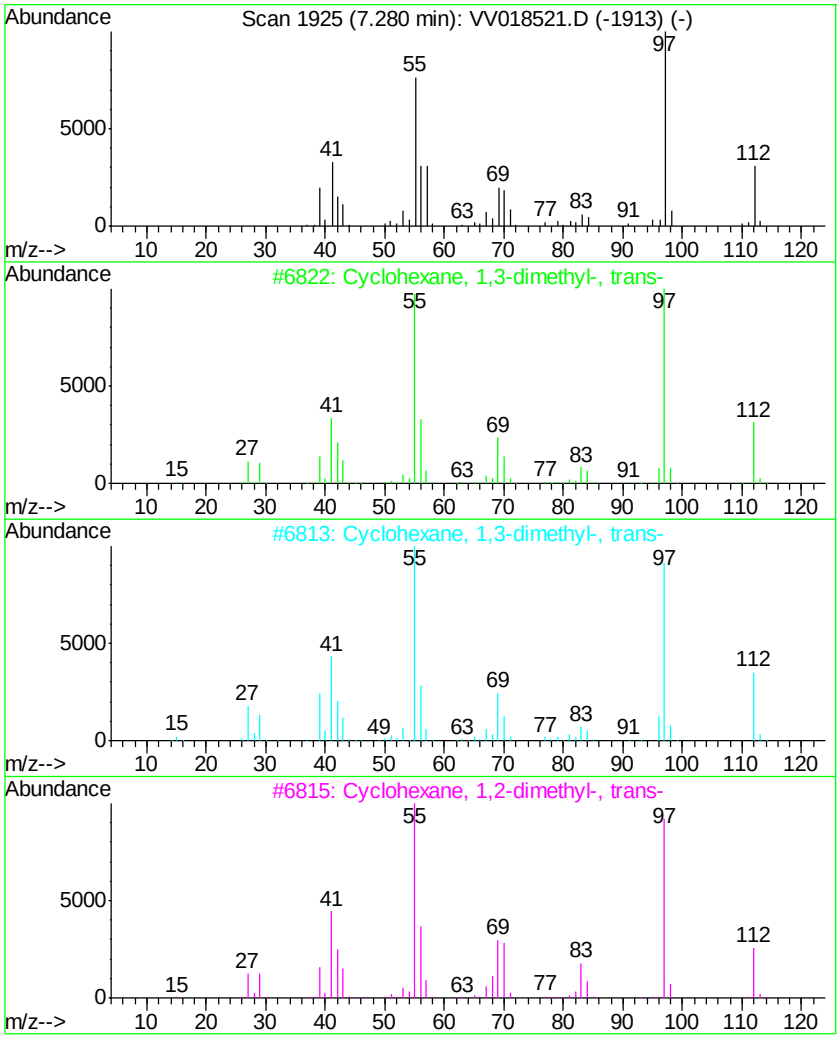
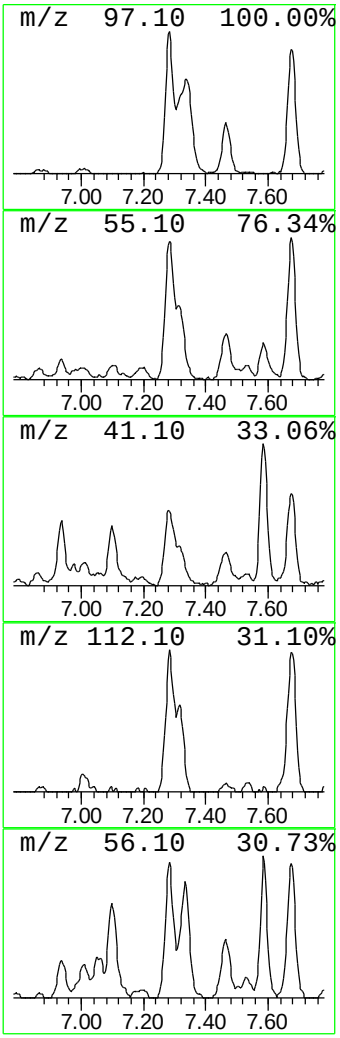
Peak Number 12 (DEL) Alkane: Cyclic7.28 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.28	6.20 ug/L	143522	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
2	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	90
3	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	87
4	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	87
5	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	86





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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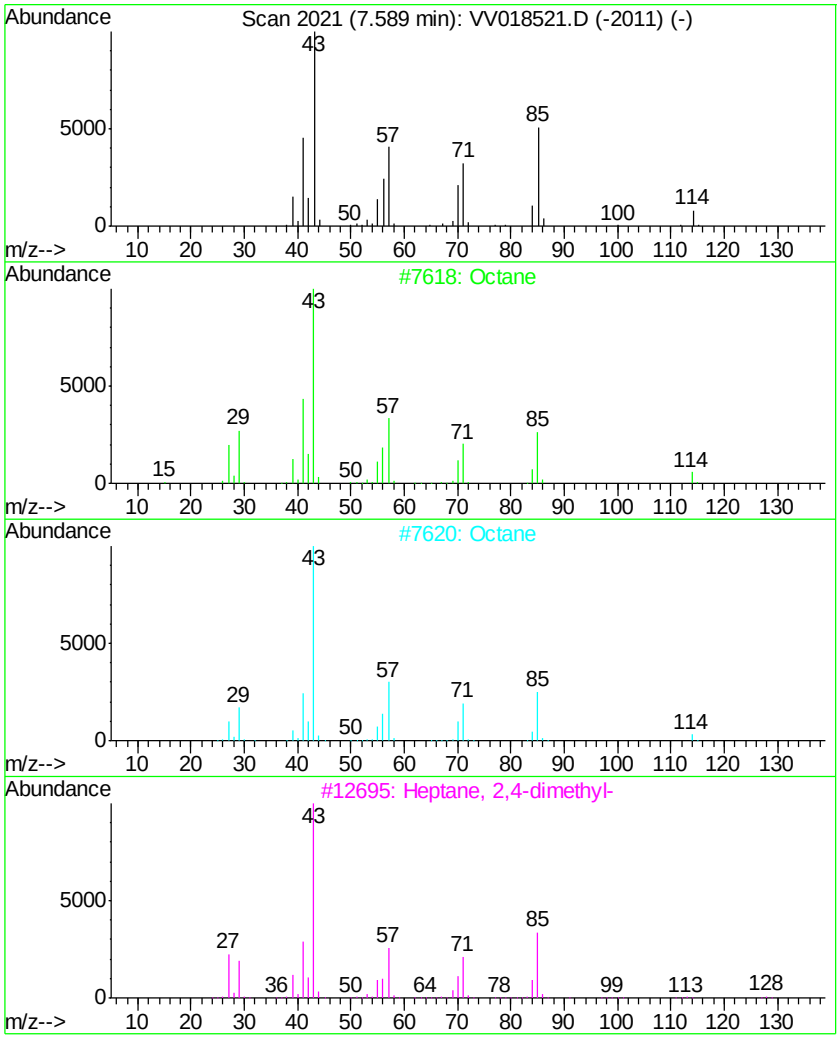
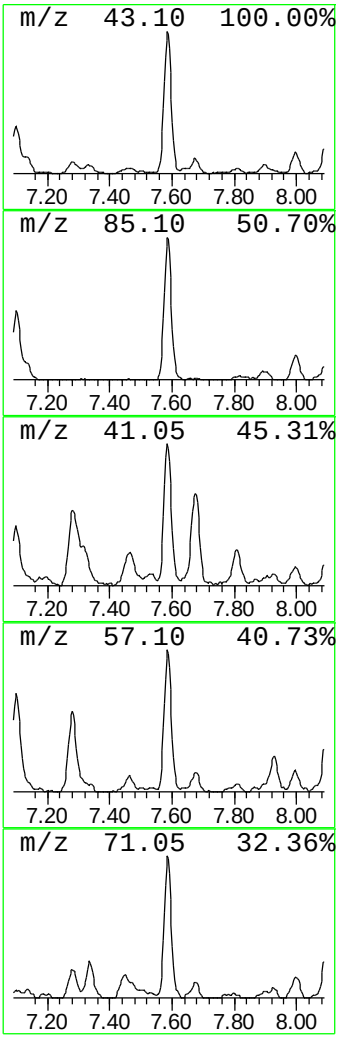
Peak Number 13 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.59	5.65 ug/L	130791	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane	114	C8H18	000111-65-9	94
2	Octane	114	C8H18	000111-65-9	91
3	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	78
4	Oxalic acid, isohexyl pentyl ester	244	C13H24O4	1000309-32-8	78
5	Heptane, 2,4-dimethyl-	128	C9H20	002213-23-2	72





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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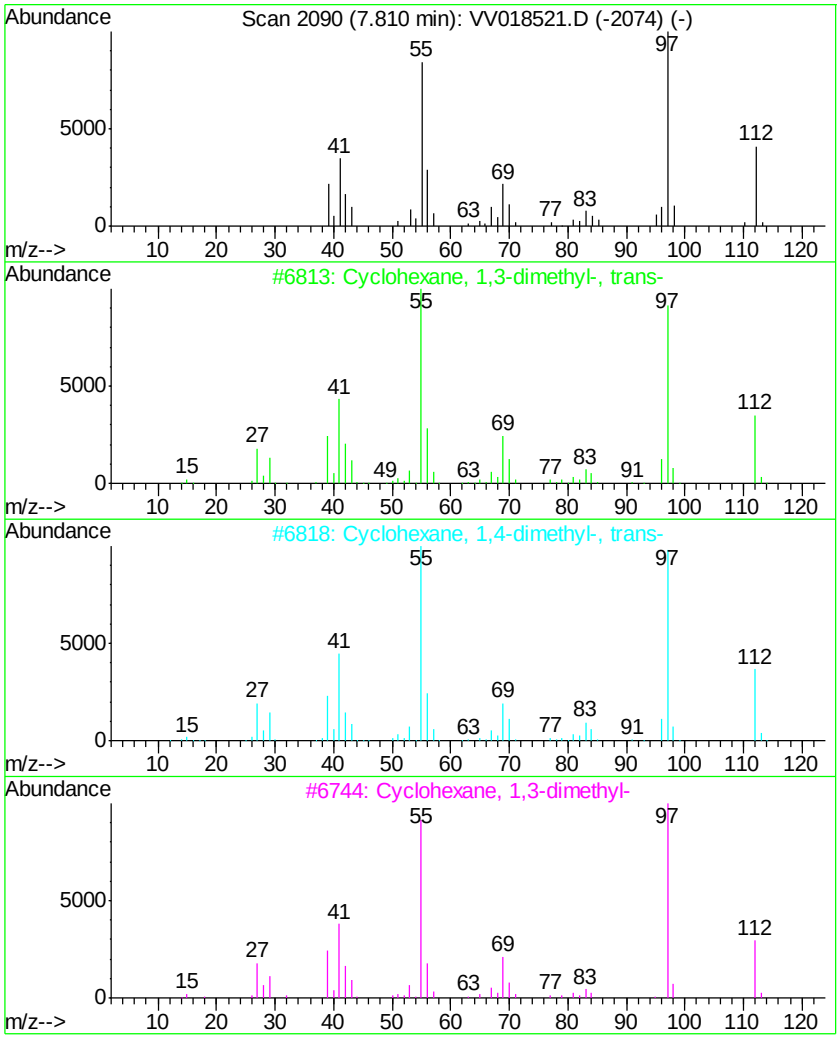
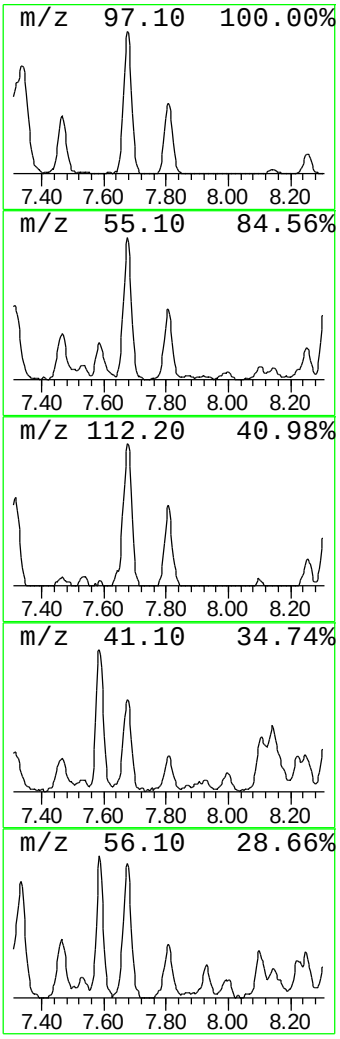
Peak Number 14 (DEL) Alkane: Cyclic7.81 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.81	2.67 ug/L	61745	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	96
2	Cyclohexane, 1,4-dimethyl-, trans-	112	C8H16	002207-04-7	94
3	Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	93
4	Cyclohexane, 1,3-dimethyl-, trans-	112	C8H16	002207-03-6	91
5	Cyclohexane, 1,4-dimethyl-	112	C8H16	000589-90-2	90





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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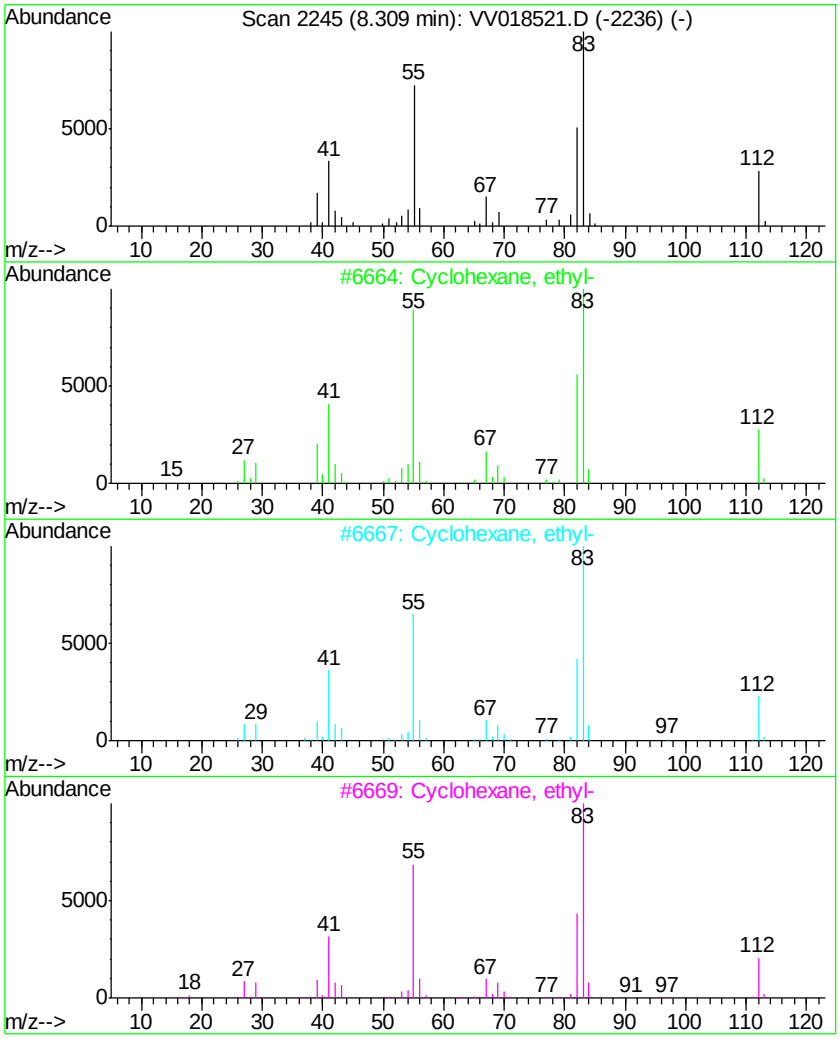
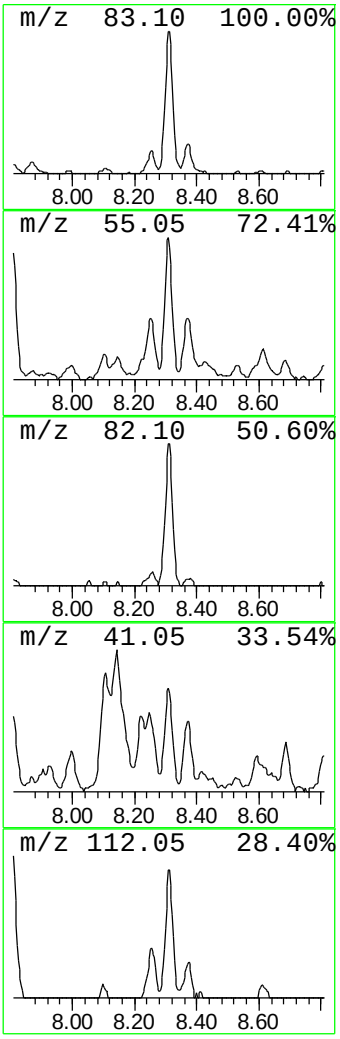
Peak Number 15 (DEL) Alkane: Cyclic8.31 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.31	2.51 ug/L	58168	Chlorobenzene-d5	8.87

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, ethyl-	112	C8H16	001678-91-7	95
2	Cyclohexane, ethyl-	112	C8H16	001678-91-7	91
3	Cyclohexane, ethyl-	112	C8H16	001678-91-7	91
4	Cyclohexane, ethyl-	112	C8H16	001678-91-7	91
5	2-Hexene, 2,3-dimethyl-	112	C8H16	007145-20-2	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

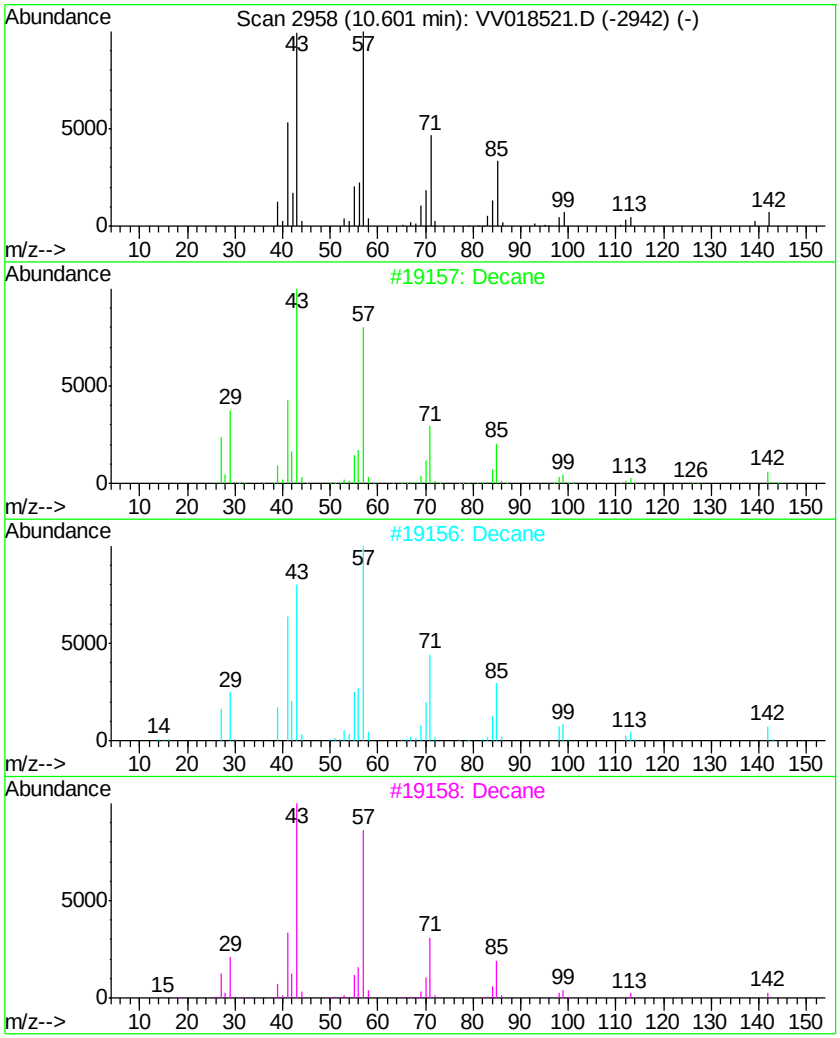
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

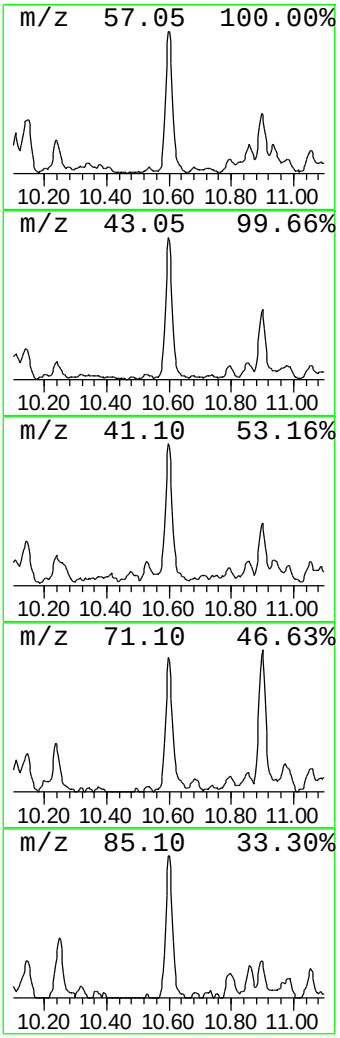
****

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.60	5.50 ug/L	149396	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	95
2	Decane	142	C10H22	000124-18-5	94
3	Decane	142	C10H22	000124-18-5	91
4	Undecane, 2,7-dimethyl-	184	C13H28	017301-24-5	80
5	Decane	142	C10H22	000124-18-5	76





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

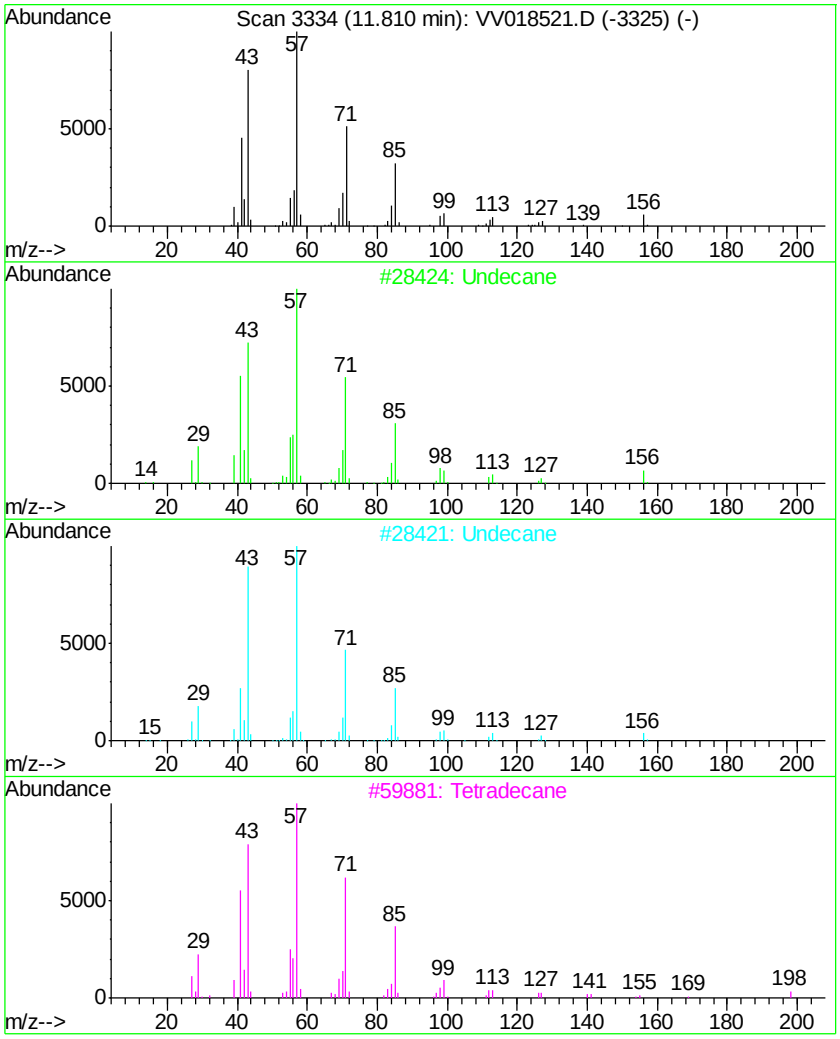
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVLM092920WMA.M
Quant Title : VOC Analysis

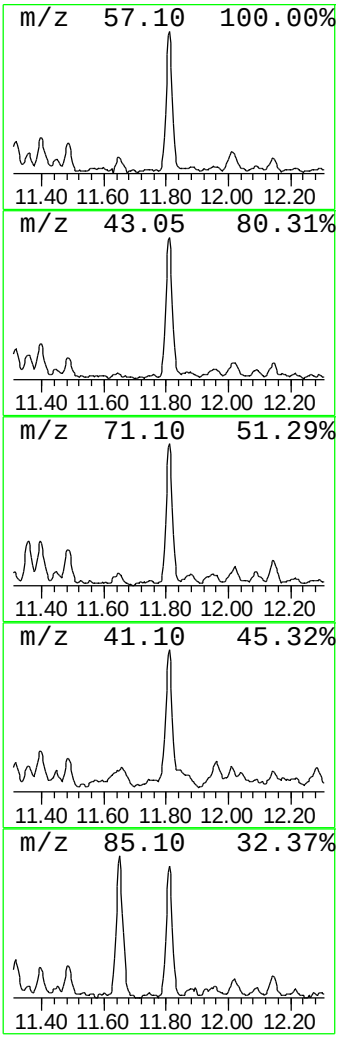
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.81	4.78 ug/L	129799	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane	156	C11H24	001120-21-4	96
2	Undecane	156	C11H24	001120-21-4	93
3	Tetradecane	198	C14H30	000629-59-4	90
4	Decane	142	C10H22	000124-18-5	80
5	Decane	142	C10H22	000124-18-5	80





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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

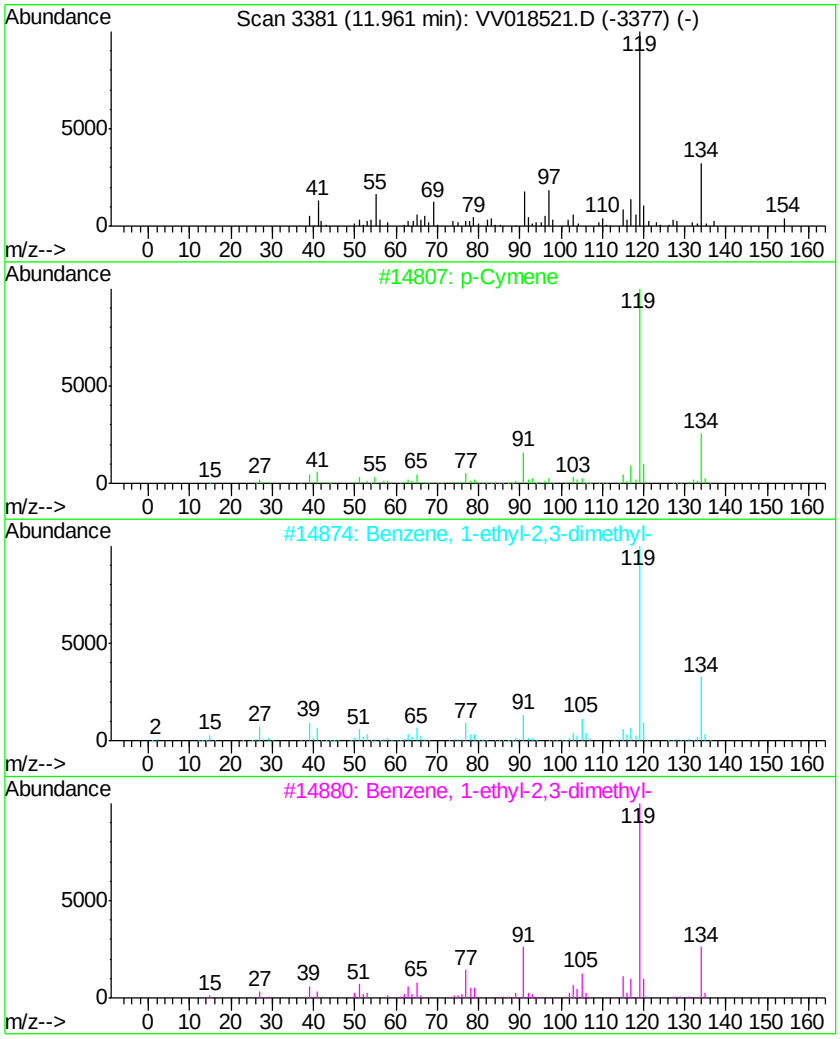
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

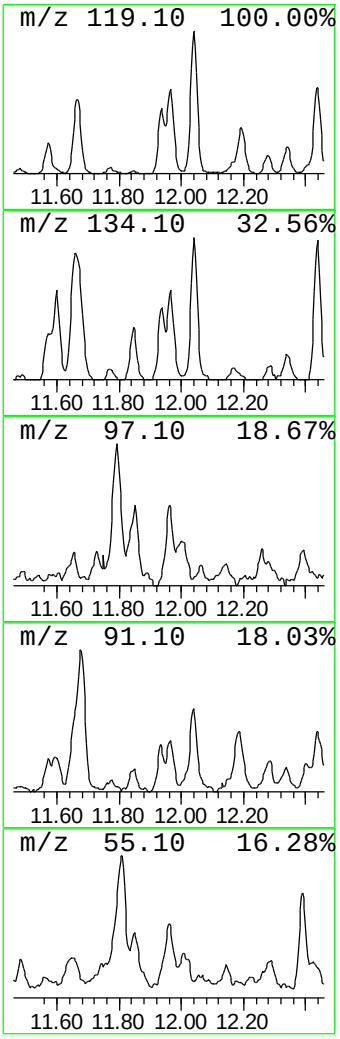
****

Peak Number 18 p-Cymene Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.96	2.63 ug/L	71588	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	87
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	83
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	83
4	o-Cymene	134	C10H14	000527-84-4	83
5	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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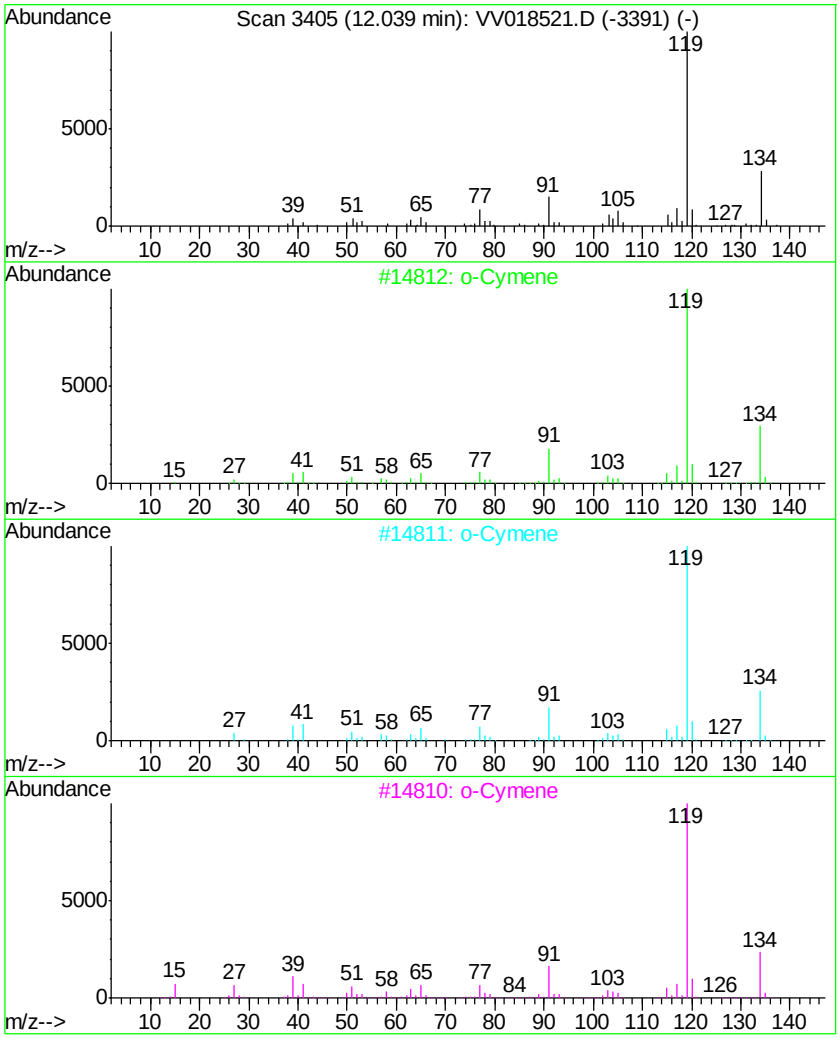
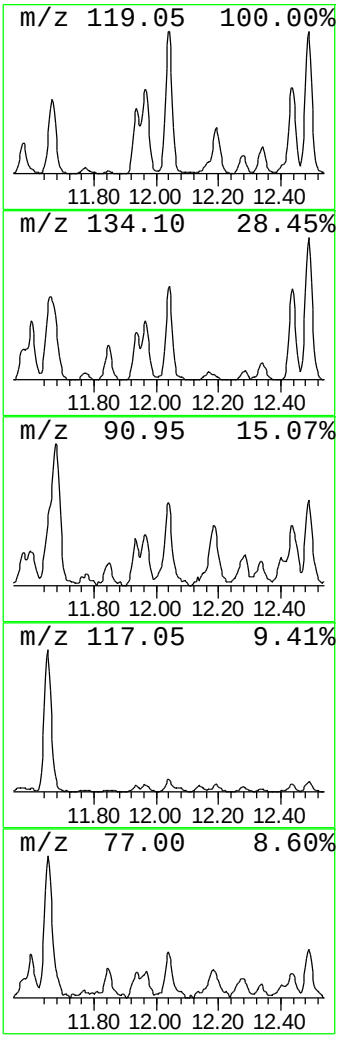
Peak Number 19 o-Cymene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.04	3.94 ug/L	107146	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	97
2	o-Cymene	134	C10H14	000527-84-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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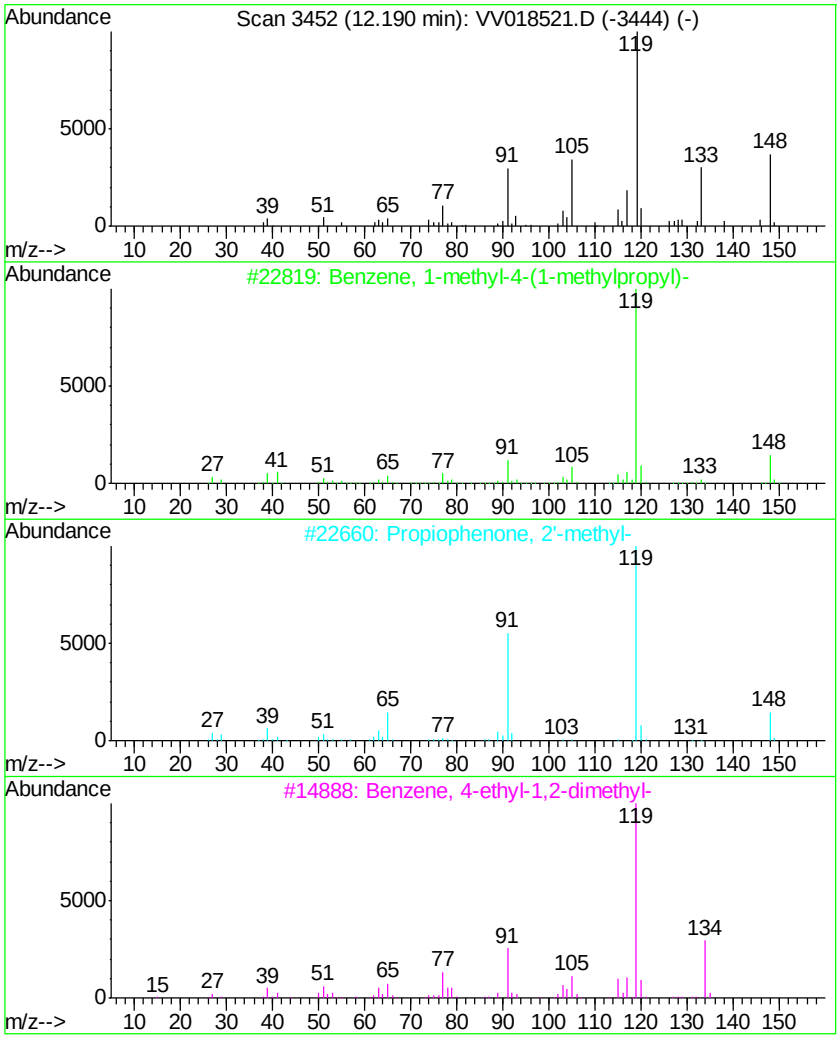
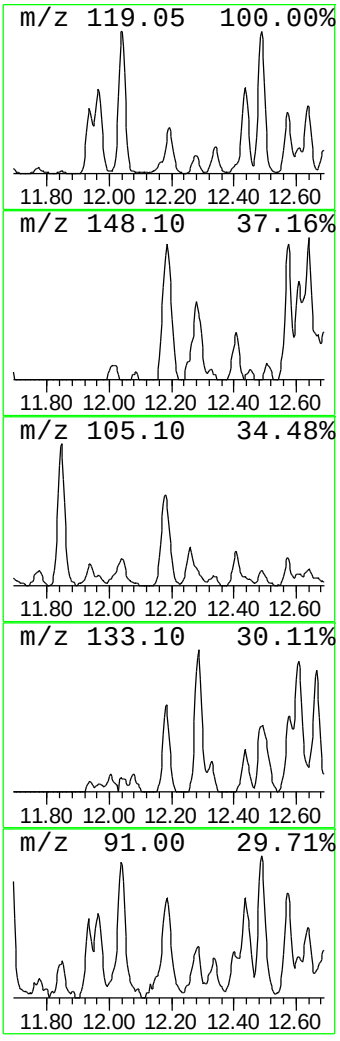
Peak Number 20 Benzene, 1-methyl-4-(1-meth... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.19	2.80 ug/L	76125	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	58
2	Propiophenone, 2'-methyl-	148	C10H12O	002040-14-4	53
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	38
4	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	38
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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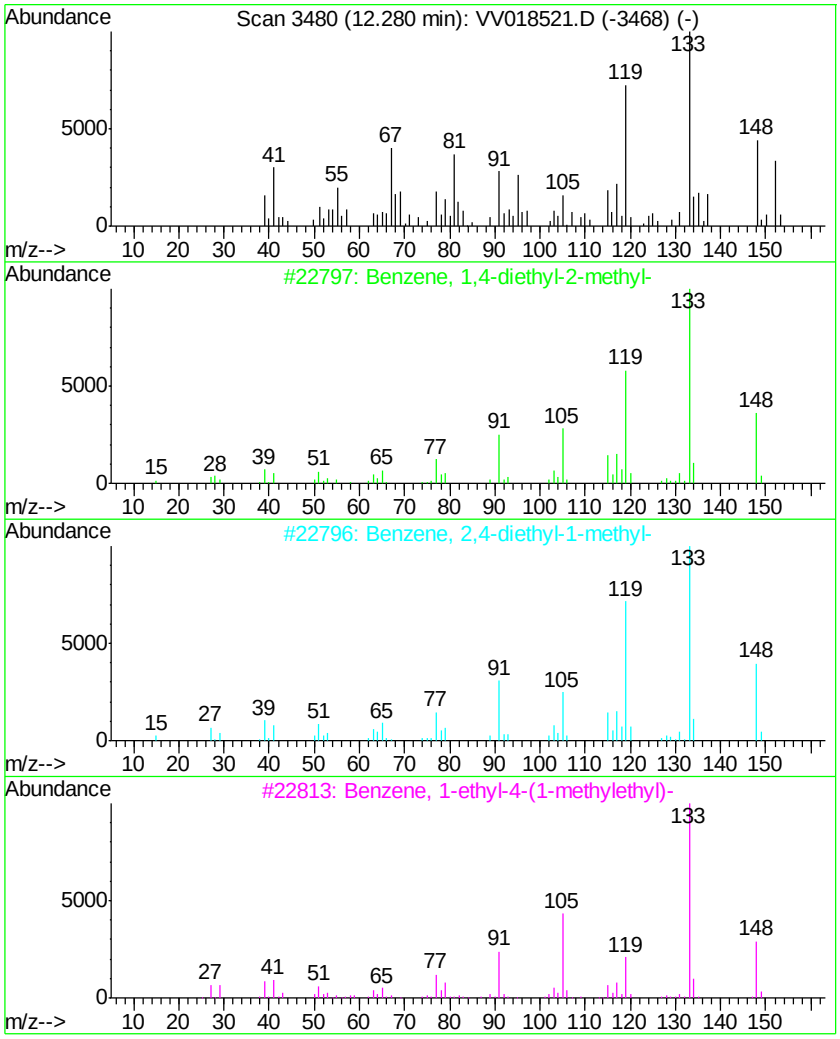
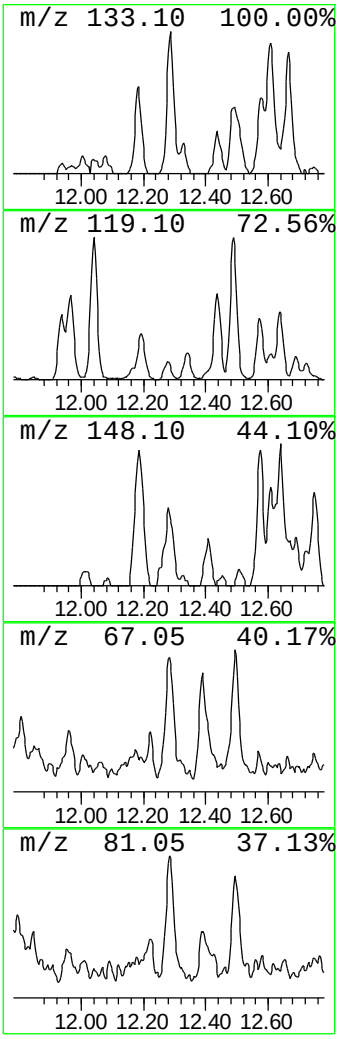
Peak Number 21 Benzene, 1,4-diethyl-2-methyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.28	2.59 ug/L	70262	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	64
2	Benzene, 2,4-diethyl-1-methyl-	148	C11H16	001758-85-6	58
3	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	49
4	Benzene, pentamethyl-	148	C11H16	000700-12-9	46
5	1,5,6,7-Tetramethylbicyclo[3.2.0...	148	C11H16	134329-46-7	43





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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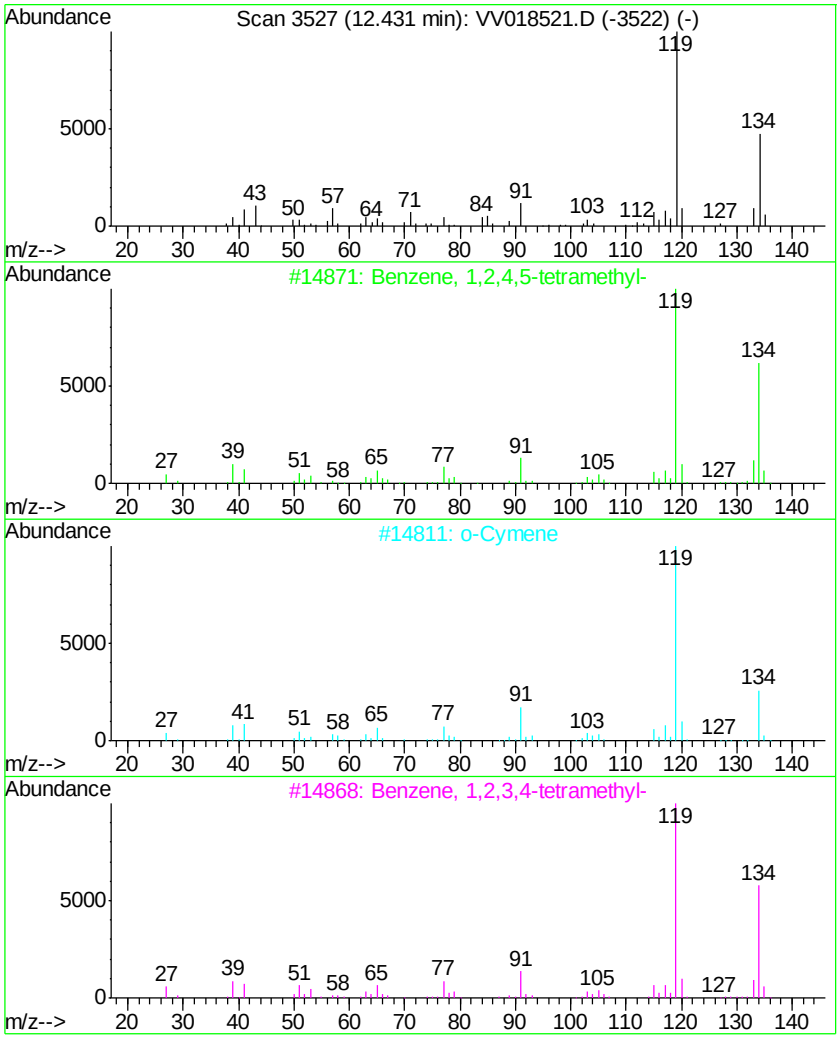
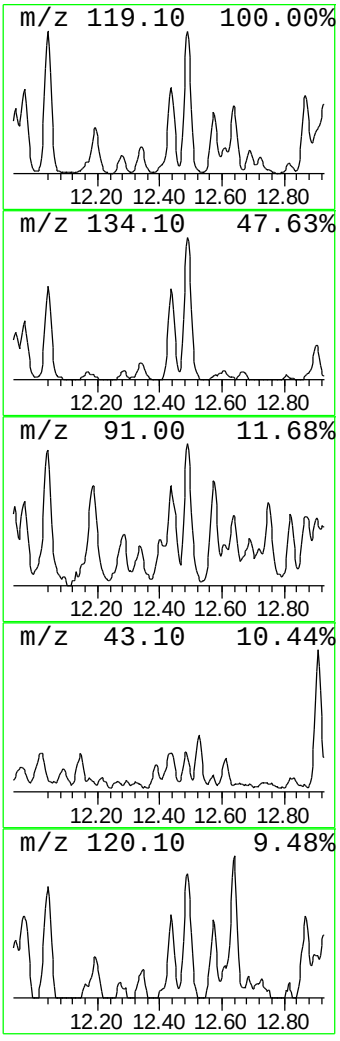
Peak Number 22 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.43	2.65 ug/L	71928	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2	o-Cymene	134	C10H14	000527-84-4	93
3	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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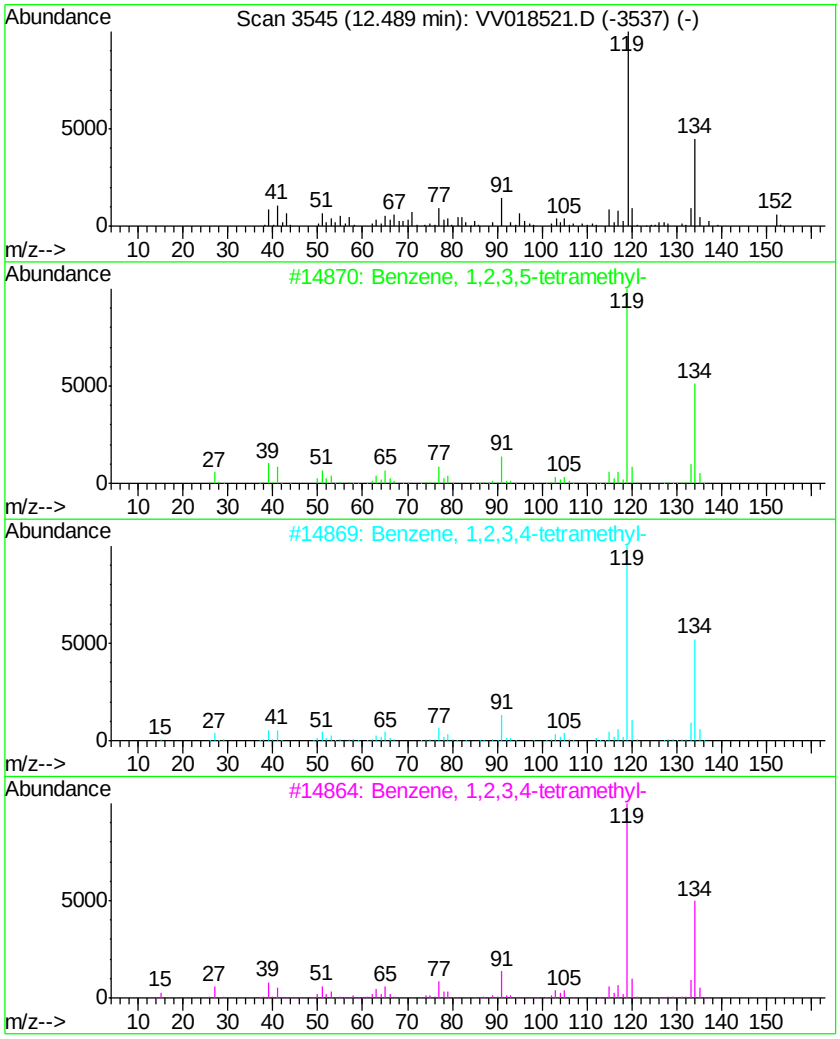
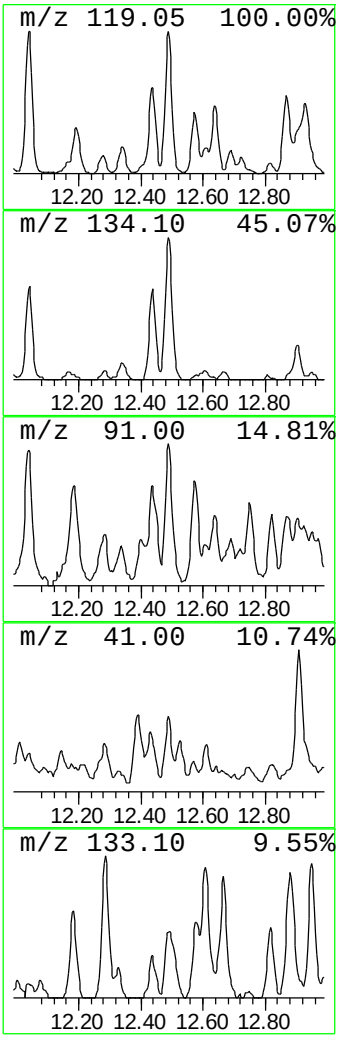
Peak Number 23 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.49	4.50 ug/L	122171	1,4-Dichlorobenzene-d4	11.28

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





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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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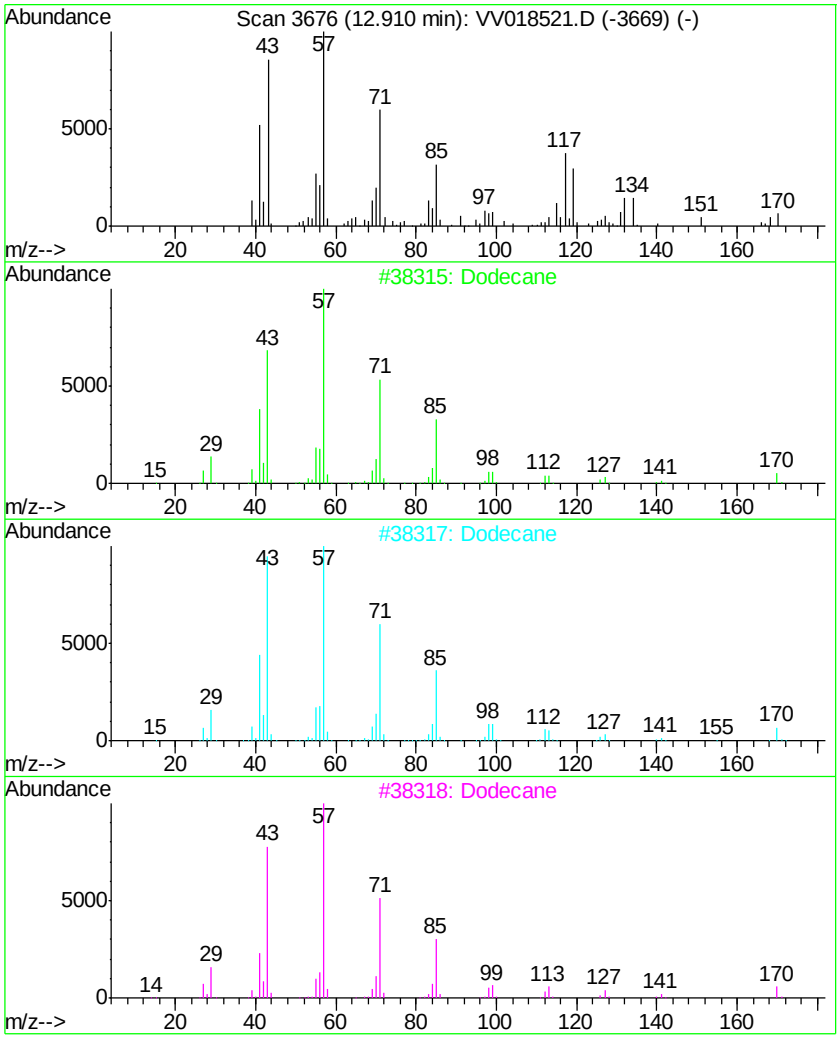
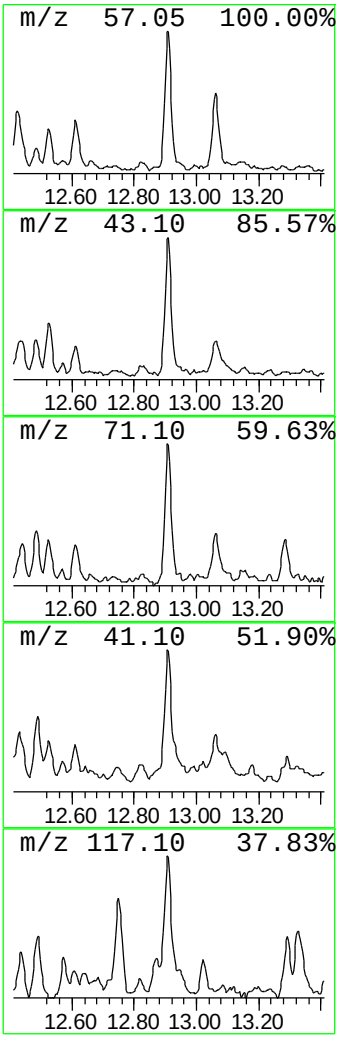
Peak Number 24 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.91	6.67 ug/L	181213	1,4-Dichlorobenzene-d4	11.28

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane	170	C12H26	000112-40-3	64
2	Dodecane	170	C12H26	000112-40-3	50
3	Dodecane	170	C12H26	000112-40-3	50
4	Dodecane	170	C12H26	000112-40-3	50
5	Nonadecane	268	C19H40	000629-92-5	43





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV092920\
Data File : VV018521.D
Acq On : 29 Sep 2020 17:06
Operator : SY/MD
Sample : L4109-19ME
Misc : 6.45g/5.0mL/100uL/5.0mL/MSVOA_V/MEOH
ALS Vial : 16 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : BG3Z0ME

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...	2.53	206.2	ug/L	3737740	1	5.64	906572	50.0
(DEL) Alkane: Str...	2.76	303.7	ug/L	5506780	1	5.64	906572	50.0
(DEL) Alkane: Str...	3.05	1026.2	ug/L	18605600	1	5.64	906572	50.0
(DEL) Alkane: Str...	3.55	80.2	ug/L	1454430	1	5.64	906572	50.0
(DEL) Alkane: Str...	3.68	68.5	ug/L	1242890	1	5.64	906572	50.0
(DEL) Alkane: Cyc...	3.77	65.7	ug/L	1190740	1	5.64	906572	50.0
(DEL) Alkane: Str...	4.78	2.9	ug/L	51998	1	5.64	906572	50.0
(DEL) Alkane: Str...	4.93	3.0	ug/L	54835	1	5.64	906572	50.0
(DEL) Alkane: Str...	6.94	3.4	ug/L	61942	1	5.64	906572	50.0
(DEL) Alkane: Str...	7.10	3.8	ug/L	69260	1	5.64	906572	50.0
(DEL) Alkane: Cyc...	7.28	6.2	ug/L	143522	2	8.87	1157820	50.0
(DEL) Alkane: Str...	7.59	5.7	ug/L	130791	2	8.87	1157820	50.0
(DEL) Alkane: Cyc...	7.81	2.7	ug/L	61745	2	8.87	1157820	50.0
(DEL) Alkane: Cyc...	8.31	2.5	ug/L	58168	2	8.87	1157820	50.0
(DEL) Alkane: Str...	10.60	5.5	ug/L	149396	3	11.28	1358570	50.0
(DEL) Alkane: Str...	11.81	4.8	ug/L	129799	3	11.28	1358570	50.0
p-Cymene	11.96	2.6	ug/L	71588	3	11.28	1358570	50.0
o-Cymene	12.04	3.9	ug/L	107146	3	11.28	1358570	50.0
Benzene, 1-methyl...	12.19	2.8	ug/L	76125	3	11.28	1358570	50.0
Benzene, 1,4-diet...	12.28	2.6	ug/L	70262	3	11.28	1358570	50.0
Benzene, 1,2,4,5-...	12.43	2.6	ug/L	71928	3	11.28	1358570	50.0
Benzene, 1,2,3,5-...	12.49	4.5	ug/L	122171	3	11.28	1358570	50.0
(DEL) Alkane: Str...	12.91	6.7	ug/L	181213	3	11.28	1358570	50.0