LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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\SOMVTR112219WMA.M
Title : TRACE VOA SOM01.0

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.209	28	37	62	rBV	647486	1274588	74.08%	6.699%
2	1.315	64	70	79	rVB	208194	234628	13.64%	1.233%
3	1.579	146	152	163	rVB	149052	166992	9.71%	0.878%
4	1.647	165	173	184	rVB	594403	721932	41.96%	3.794%
5	1.820	220	227	244	rVB	153082	207879	12.08%	1.093%
6	1.917	251	257	264	rVB	33403	41769	2.43%	0.220%
7	2.042	289	296	307	rVB2	37337	53499	3.11%	0.281%
8	2.126	313	322	345	rVB	296888	519157	30.17%	2.728%
9	2.528	434	447	450	rBV4	52186	94587	5.50%	0.497%
10	2.605	463	471	489	rVB2	49036	100699	5.85%	0.529%
11	2.791	517	529	540	rBV	58679	110285	6.41%	0.580%
12	2.975	573	586	599	rBV6	9671	24130	1.40%	0.127%
13	3.074	606	617	629	rBV3	12717	25621	1.49%	0.135%
14	3.193	642	654	660	rBV10	9444	17218	1.00%	0.090%
15	3.254	665	673	682	rVV5	13903	31818	1.85%	0.167%
16	3.309	683	690	702	rVB3	17508	35025	2.04%	0.184%
17	3.409	707	721	728	rBV7	10304	22323	1.30%	0.117%
18	3.621	777	787	799	rVB6	14331	29229	1.70%	0.154%
19	3.804	828	844	861	rBV2	78230	201286	11.70%	1.058%
20	3.939	874	886	918	rBV2	146681	486353	28.27%	2.556%
21	4.396	1010	1028	1050	rBV	210893	548866	31.90%	2.885%
22	4.720	1117	1129	1143	rBV6	40624	99944	5.81%	0.525%
23	5.090	1224	1244	1259	rBV4	481569	1202786	69.90%	6.321%
24	5.270	1297	1300	1320	rVB4	10832	25122	1.46%	0.132%
25	5.659	1407	1421	1439	rBV	415572	842515	48.97%	4.428%
26	6.113	1547	1562	1578	rBV	275444	574542	33.39%	3.020%
27	6.843	1778	1789	1798	rVB6	9535	17504	1.02%	0.092%
28	7.026	1832	1846	1863	rBV	132200	243942	14.18%	1.282%
29	7.357	1935	1949	1961	rBV	541275	935922	54.39%	4.919%
30	7.428	1961	1971	1991	rVB	303484	523947	30.45%	2.754%
31	7.659	2033	2043	2060	rBV	81886	142546	8.28%	0.749%
32	8.135	2180	2191	2224	rBV	573972	1277784	74.26%	6.716%
33	8.891	2413	2426	2444	rBV	623861	999961	58.12%	5.255%
34	9.051	2463	2476	2490	rBV	333954	536390	31.17%	2.819%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.1 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\SOMVTR112219WMA.M
Title : TRACE VOA SOM01.0

35	9.177	2501	2515	2534	rBV	1054770	1720640	100.00%	9.043%
36	9.582	2630	2641	2660	rBV	237174	375206	21.81%	1.972%
37	9.971	2753	2762	2775	rVB2	20089	33051	1.92%	0.174%
38	10.257	2838	2851	2865	rBV	229987	372770	21.66%	1.959%
39	10.392	2883	2893	2910	rBV	64934	110116	6.40%	0.579%
40	10.485	2910	2922	2941	rVV	198626	440044	25.57%	2.313%
41	10.579	2941	2951	2965	rVB2	100333	160513	9.33%	0.844%
42	10.778	3003	3013	3029	rBV	91097	150376	8.74%	0.790%
43	10.955	3056	3068	3081	rBV	400998	620938	36.09%	3.263%
44	11.289	3160	3172	3188	rBV	711349	1097015	63.76%	5.765%
45	11.376	3188	3199	3210	rVB	90228	142148	8.26%	0.747%
46	11.569	3250	3259	3270	rBV	66514	118209	6.87%	0.621%
47	11.666	3278	3289	3312	rVB	645501	1041388	60.52%	5.473%
48	11.955	3369	3379	3383	rBV2	16876	26079	1.52%	0.137%
49	12.058	3403	3411	3418	rBV2	21577	33018	1.92%	0.174%
50	12.154	3433	3441	3449	rBV	19752	33817	1.97%	0.178%
51	12.453	3525	3534	3543	rBV4	12790	23780	1.38%	0.125%
52	12.508	3543	3551	3563	rVB	19934	31040	1.80%	0.163%
53	12.768	3615	3632	3641	rBV	17953	31564	1.83%	0.166%
54	12.926	3669	3681	3694	rBV2	32659	55057	3.20%	0.289%
55	13.553	3867	3876	3891	rVB	22408	39750	2.31%	0.209%

Sum of corrected areas: 19027308

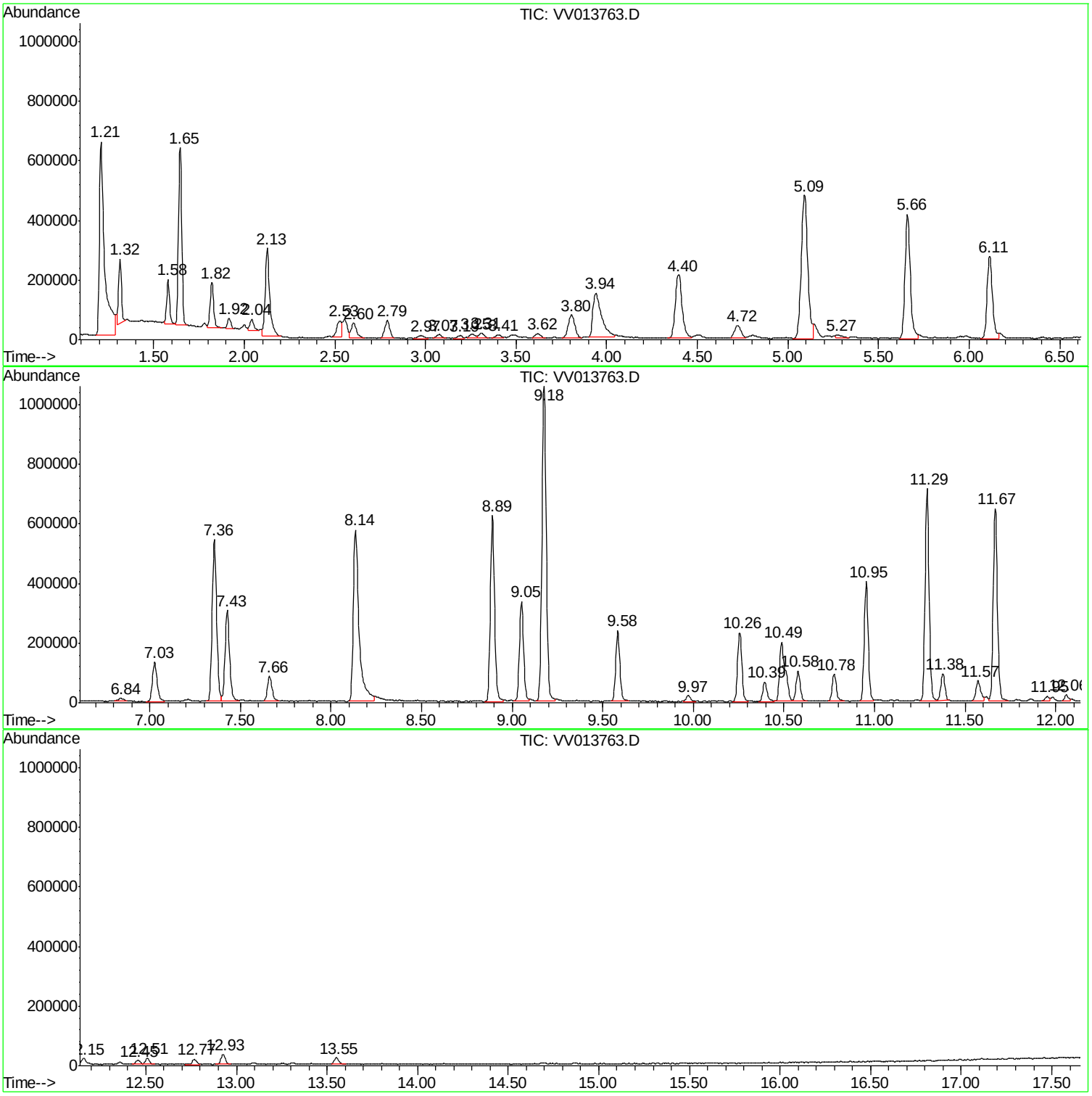
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

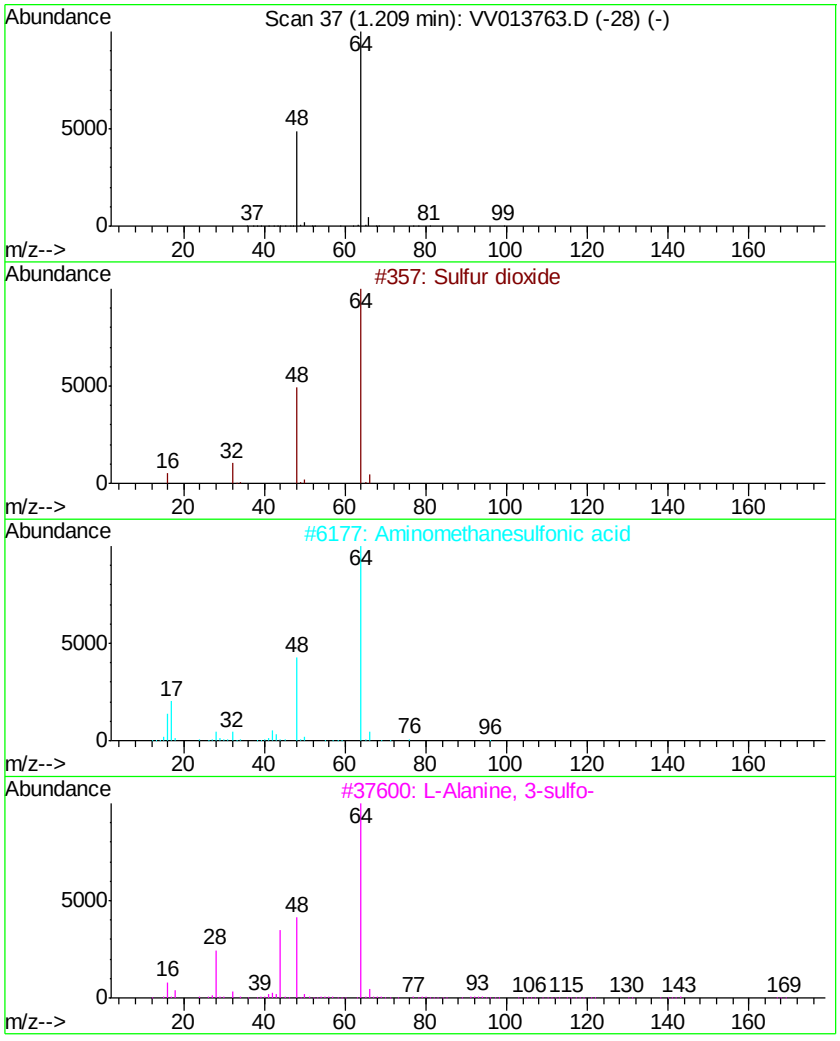
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

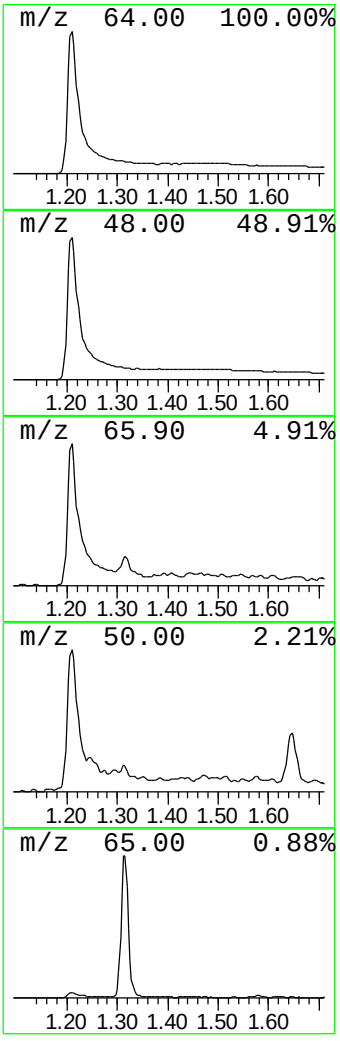
****

Peak Number 1 Sulfur dioxide Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.21	7.56 ug/L	1274590	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Sulfur dioxide	64	O2S	007446-09-5	83
2	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	74
3	L-Alanine, 3-sulfo-	169	C3H7NO5S	000498-40-8	9
4	Aminomethanesulfonic acid	111	CH5NO3S	013881-91-9	9
5	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

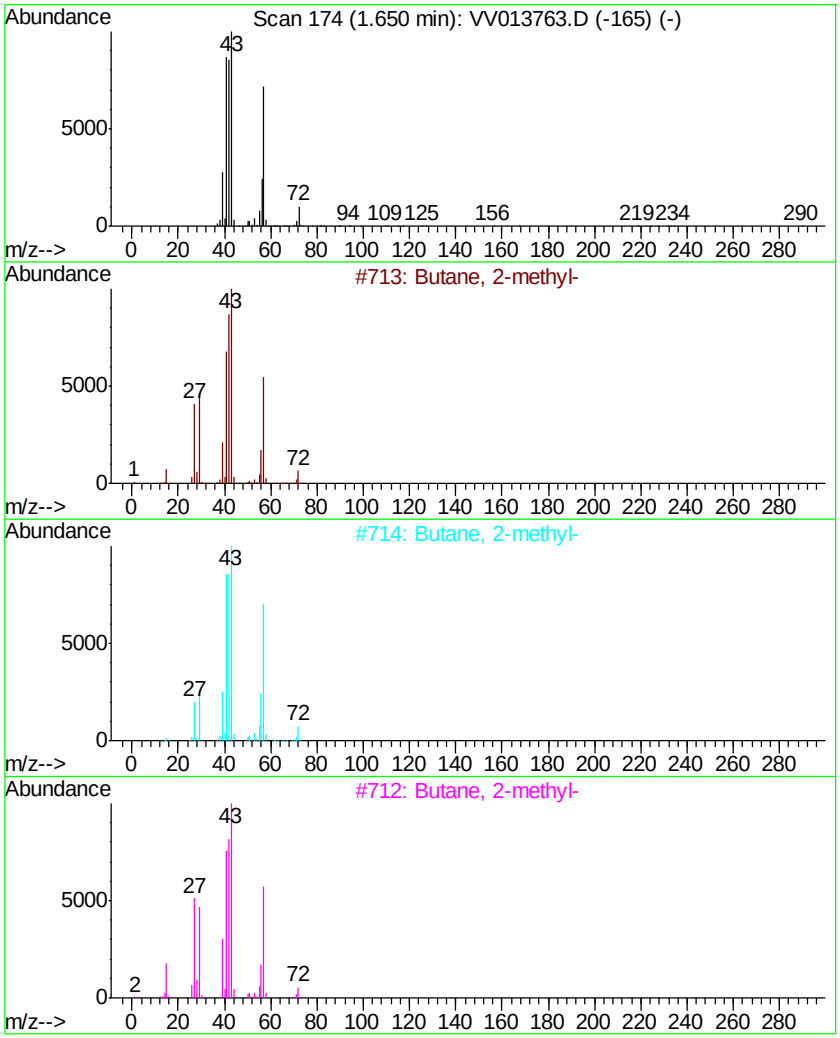
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

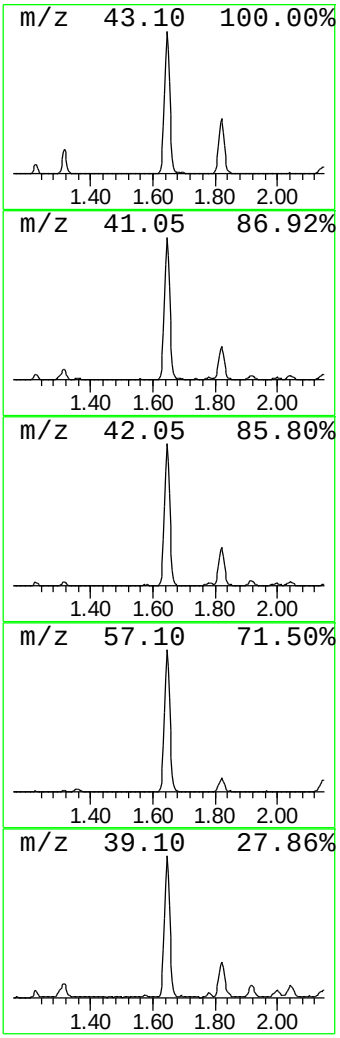
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.65	4.28 ug/L	721932	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	91
3	Butane, 2-methyl-	72	C5H12	000078-78-4	90
4	Pentane	72	C5H12	000109-66-0	45
5	Pentane	72	C5H12	000109-66-0	43





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

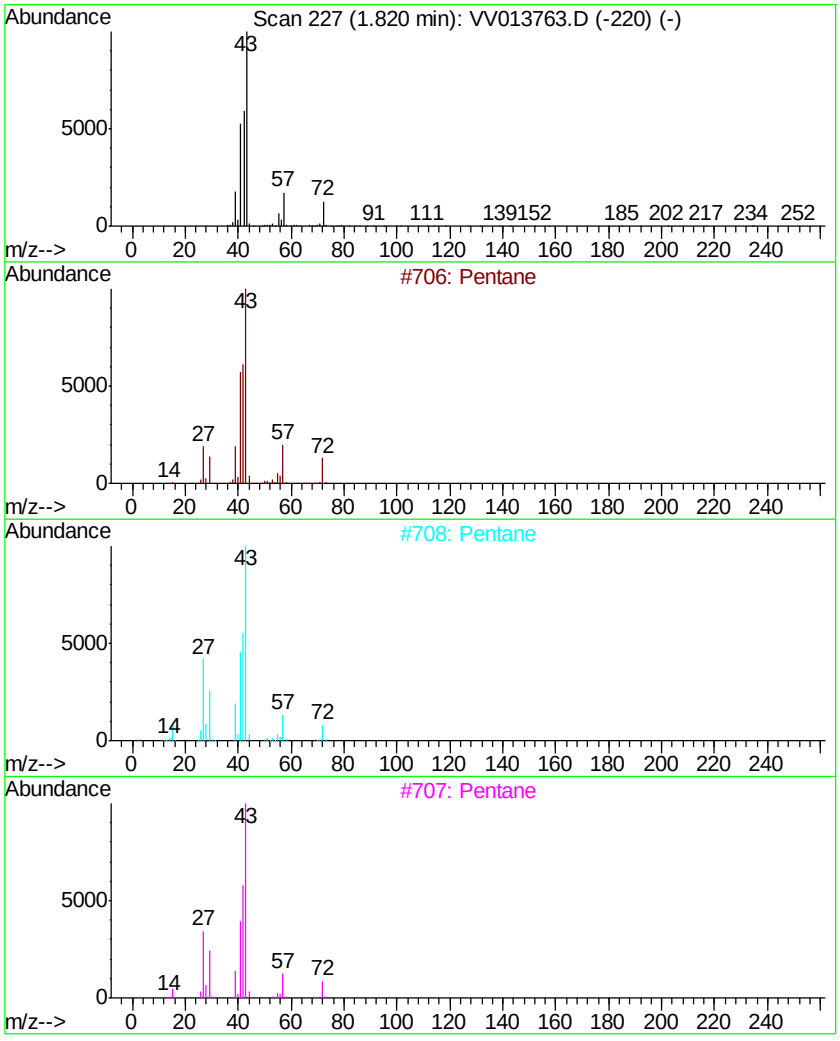
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

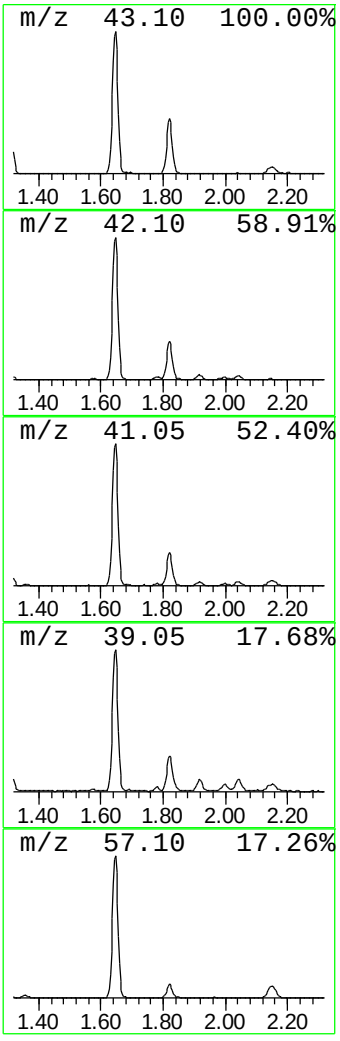
****

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.82	1.23 ug/L	207879	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	90
2	Pentane	72	C5H12	000109-66-0	86
3	Pentane	72	C5H12	000109-66-0	80
4	Pentane	72	C5H12	000109-66-0	80
5	Butane, 2-methyl-	72	C5H12	000078-78-4	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

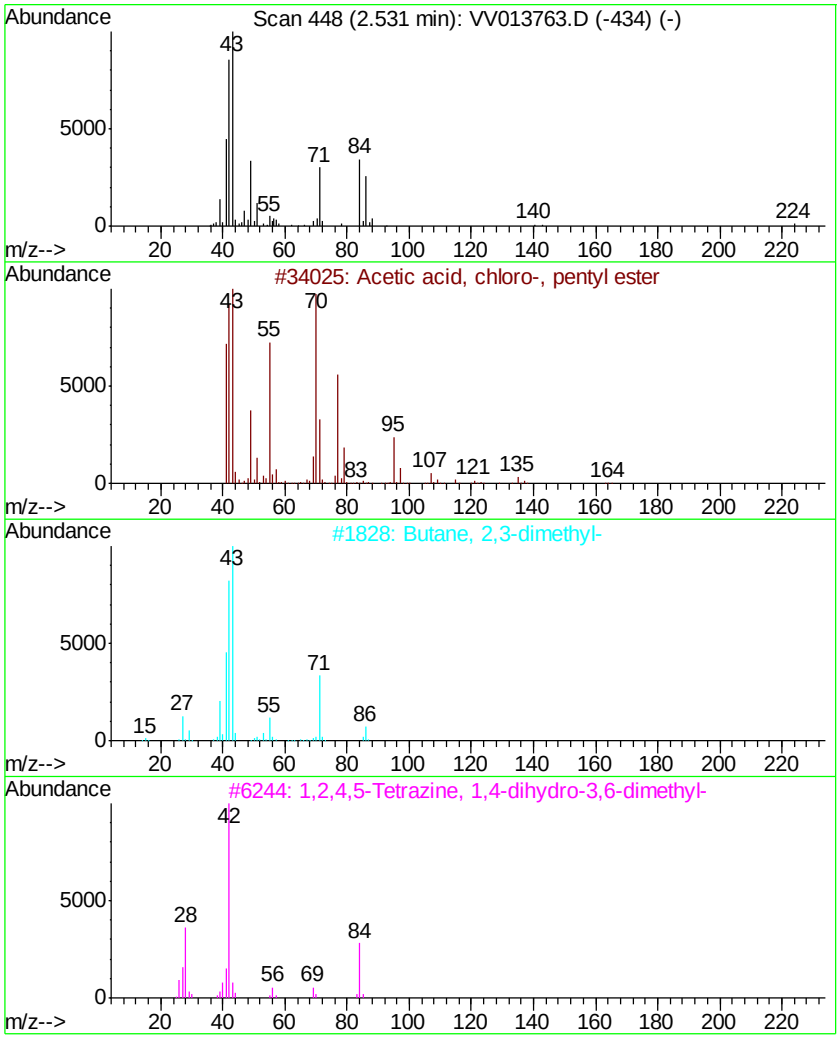
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

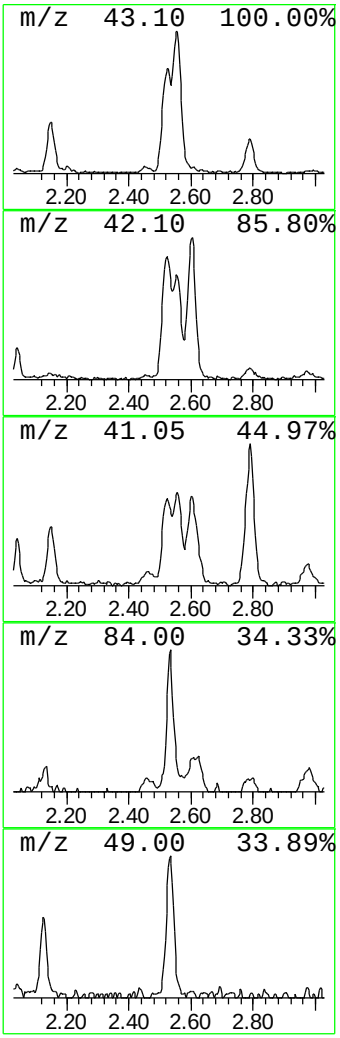
****

Peak Number 4 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.53	0.56 ug/L	94587	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acetic acid, chloro-, pentyl ester	164	C7H13ClO2	005411-55-2	38
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	35
3	1,2,4,5-Tetrazine, 1,4-dihydro-3...	112	C4H8N4	037454-64-1	27
4	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	27
5	Butyrolactone	86	C4H6O2	000096-48-0	16





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

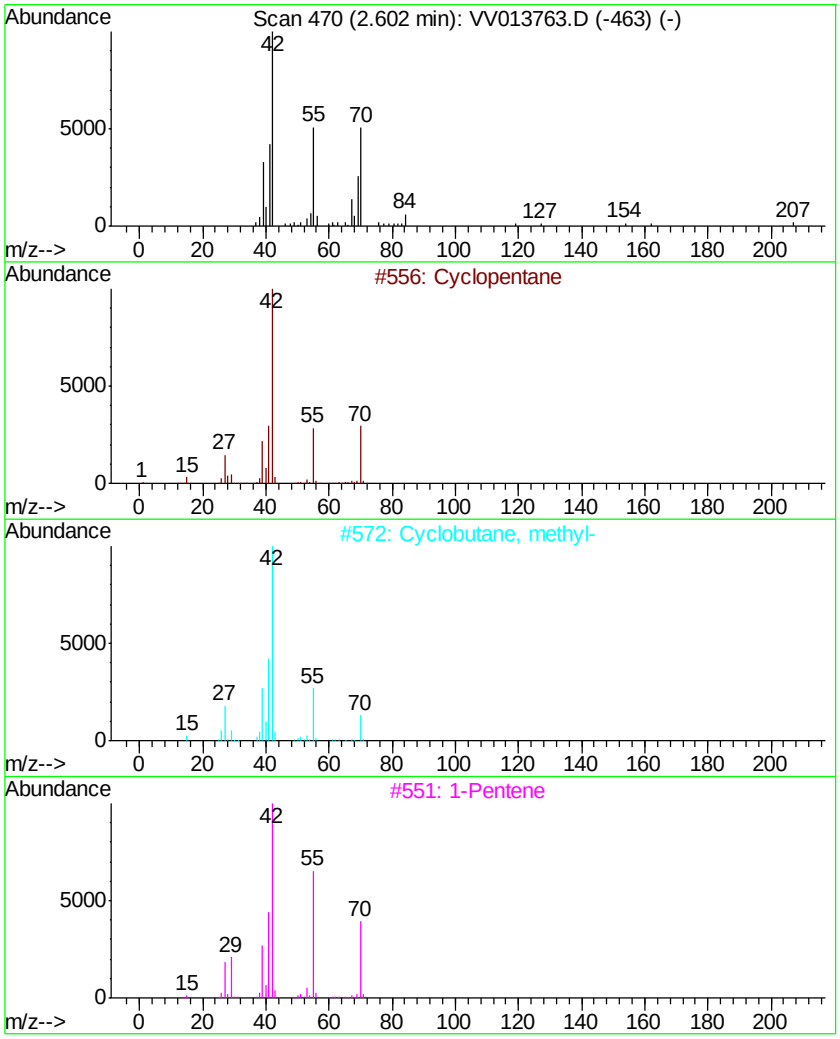
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

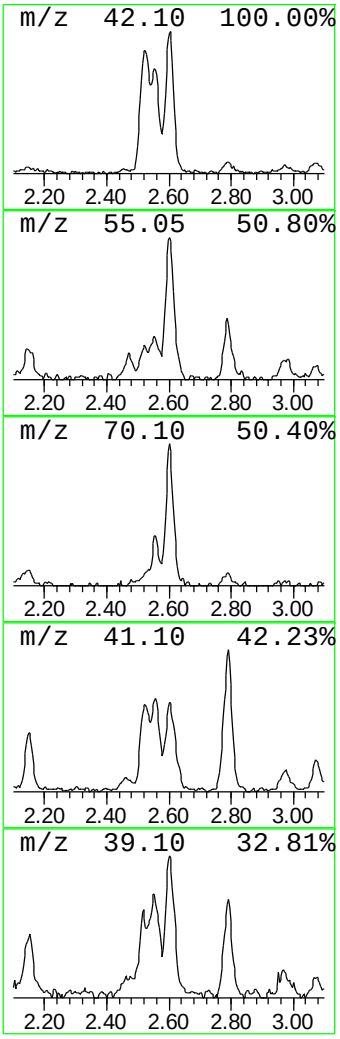
****

Peak Number 5 (DEL) Alkane: Cyclic2.60 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.60	0.60 ug/L	100699	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane	70	C5H10	000287-92-3	72
2	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
3	1-Pentene	70	C5H10	000109-67-1	72
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	59
5	Cyclopropane, ethyl-	70	C5H10	001191-96-4	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

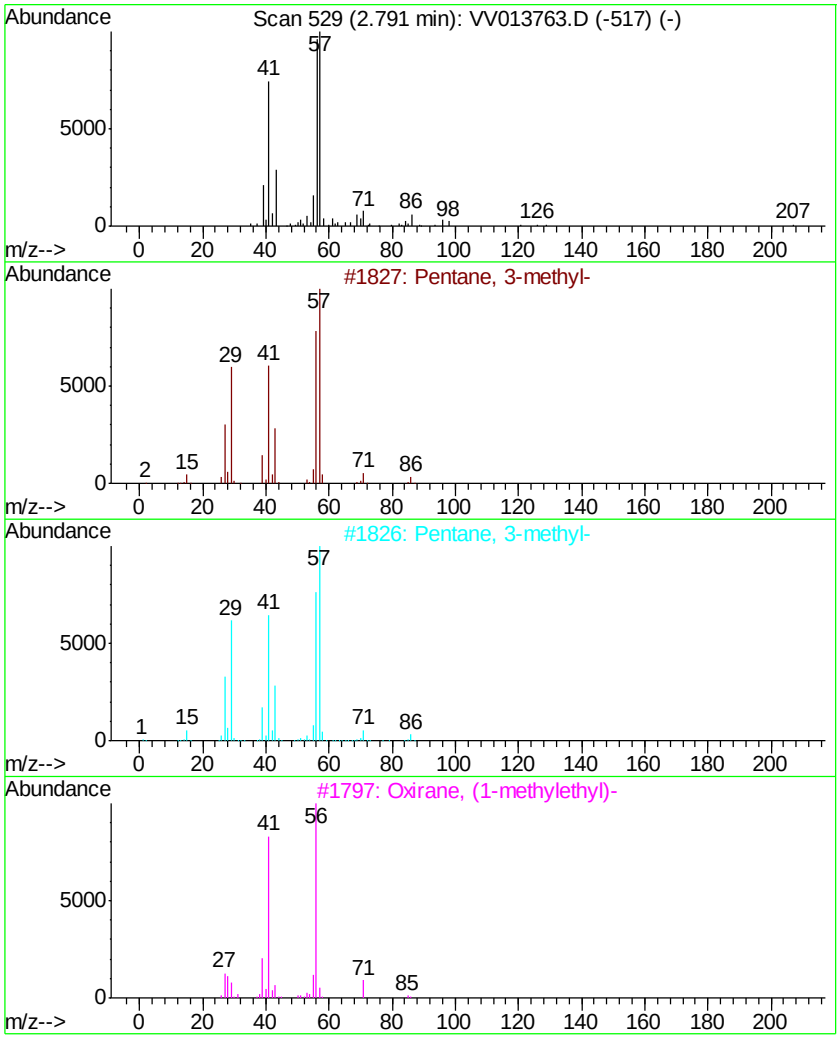
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

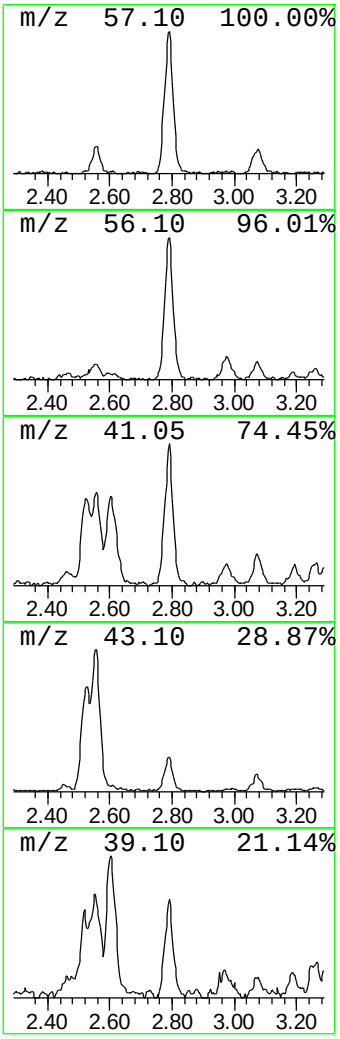
****

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.79	0.65 ug/L	110285	1,4-Difluorobenzene	5.66

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	80
2	Pentane, 3-methyl-	86	C6H14	000096-14-0	72
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	59
4	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	58
5	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	56





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

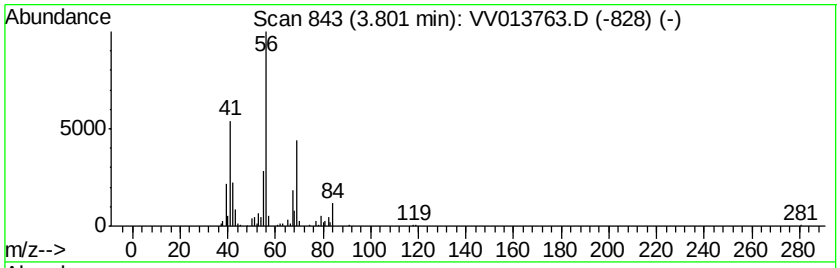
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

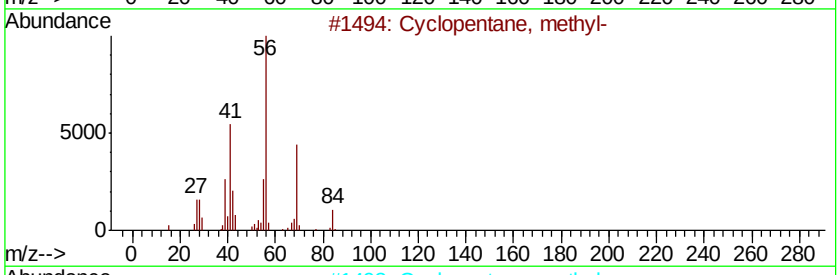
****

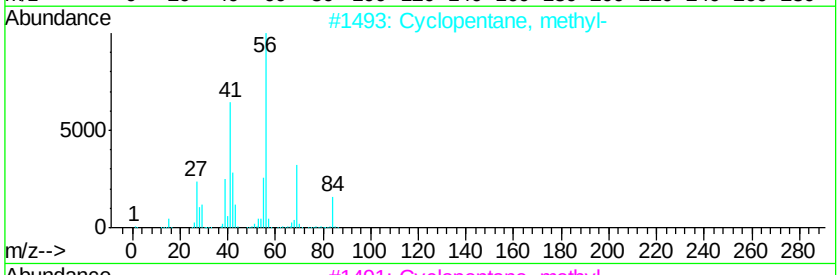
Peak Number 7 (DEL) Alkane: Cyclic3.80 Concentration Rank 7

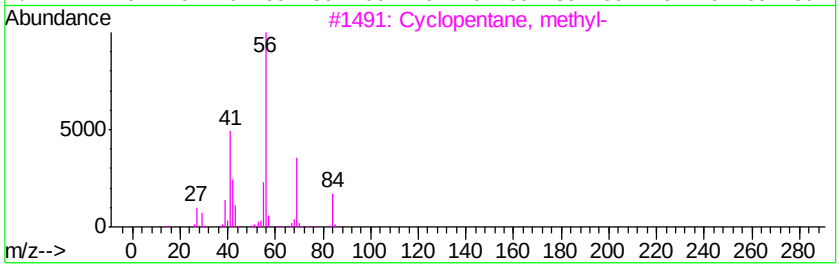
R.T.	EstConc	Area	Relative to ISTD	R.T.
3.80	1.19 ug/L	201286	1,4-Difluorobenzene	5.66

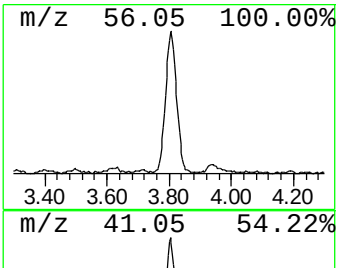
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	87
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	83
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
4	Cyclohexane	84	C6H12	000110-82-7	72
5	Cyclohexane	84	C6H12	000110-82-7	64

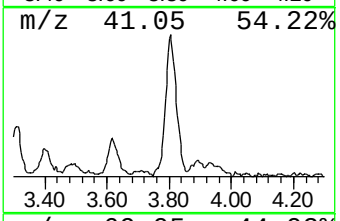


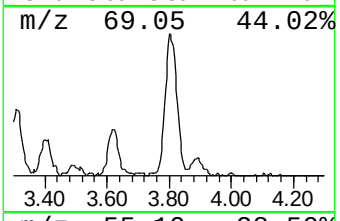


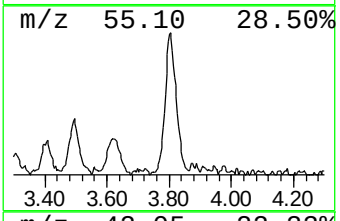


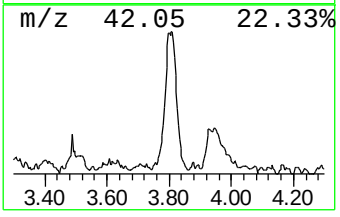












Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

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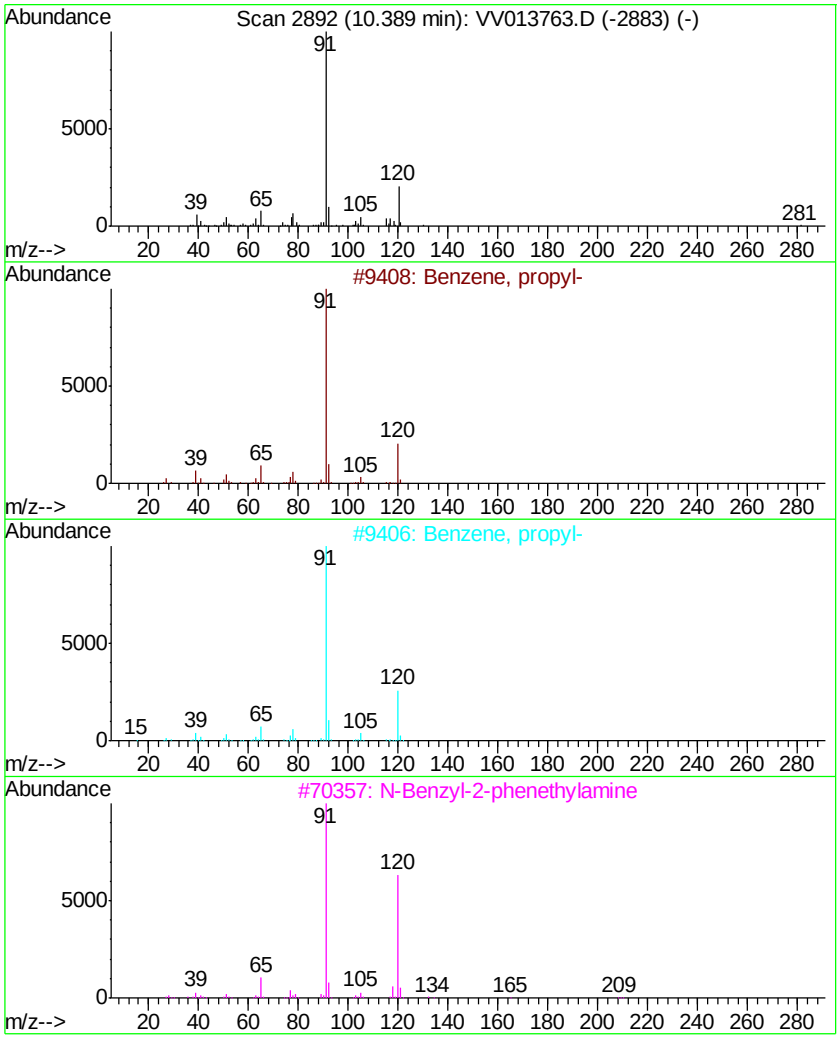
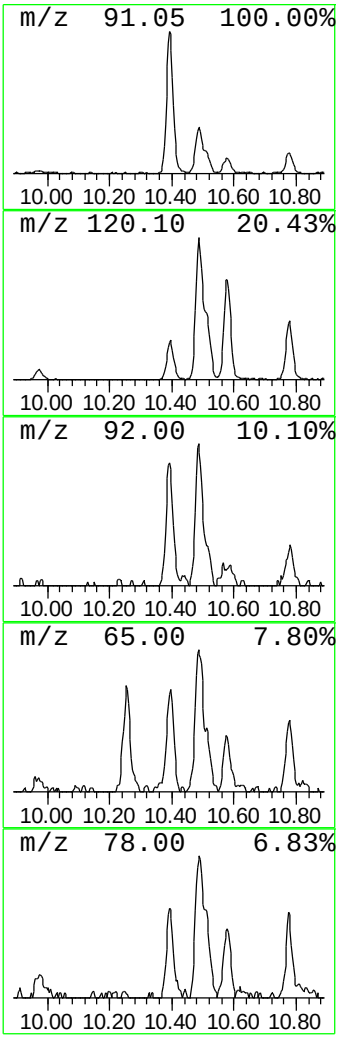
Peak Number 9 Benzene, propyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.39	0.50 ug/L	110116	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	81
2	Benzene, propyl-	120	C9H12	000103-65-1	81
3	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	64
4	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	53
5	1-Benzylamino-2-benzyloxyethane	241	C16H19NO	038336-06-0	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

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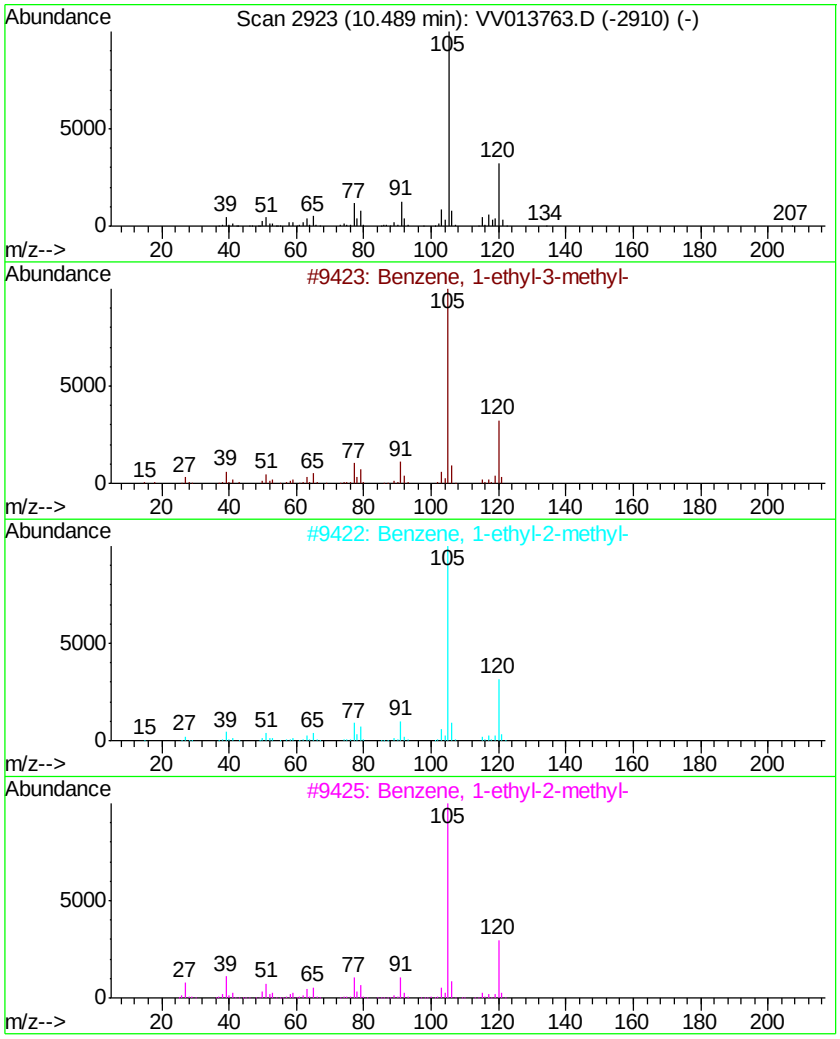
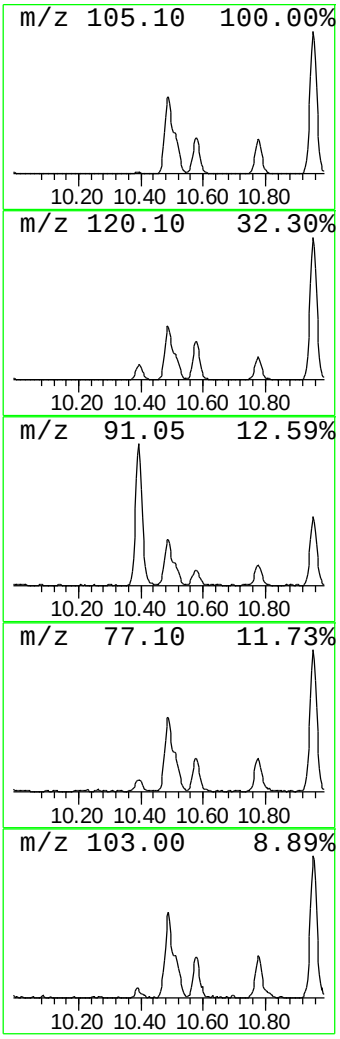
Peak Number 10 Benzene, 1-ethyl-3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	2.01 ug/L	440044	1,4-Dichlorobenzene-d4	11.29

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	94
3	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

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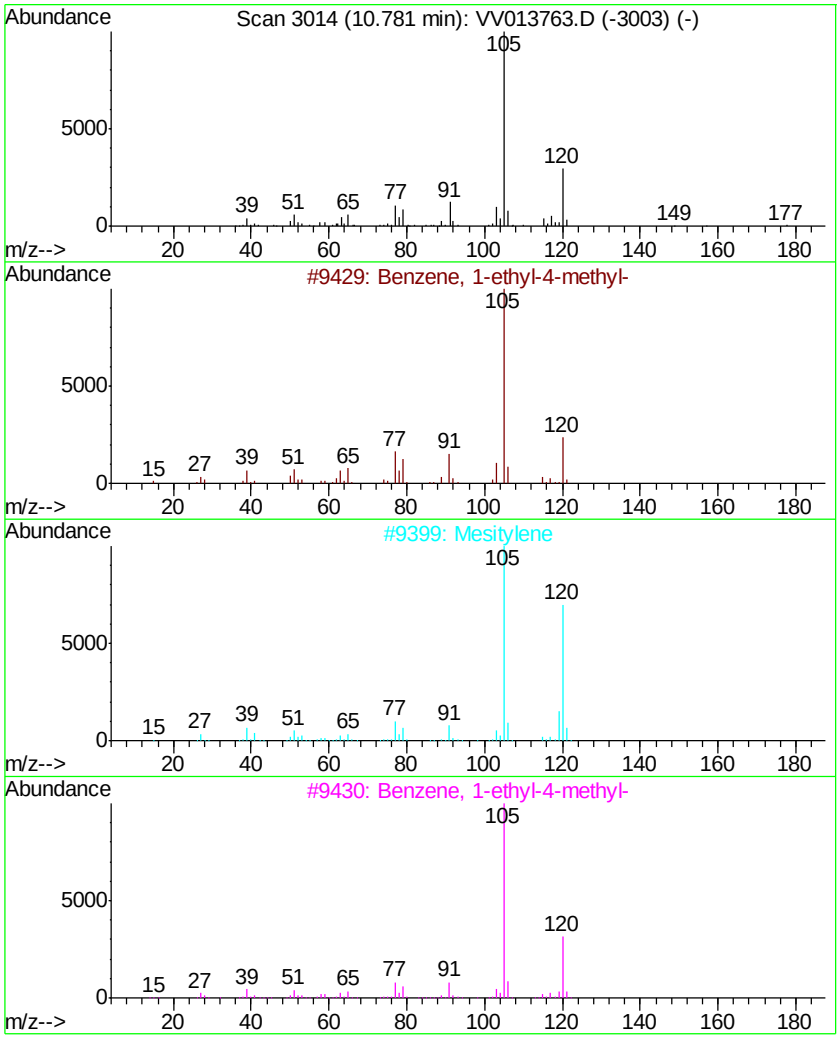
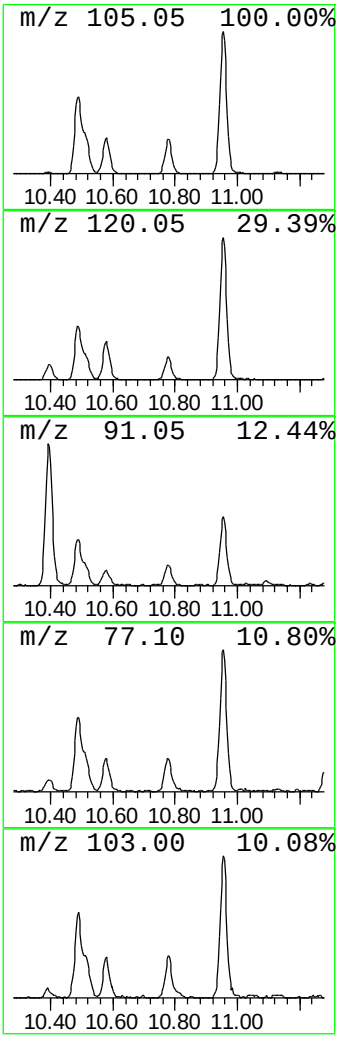
Peak Number 12 Benzene, 1-ethyl-4-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.78	0.69 ug/L	150376	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
2	Mesitylene	120	C9H12	000108-67-8	91
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5	Mesitylene	120	C9H12	000108-67-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

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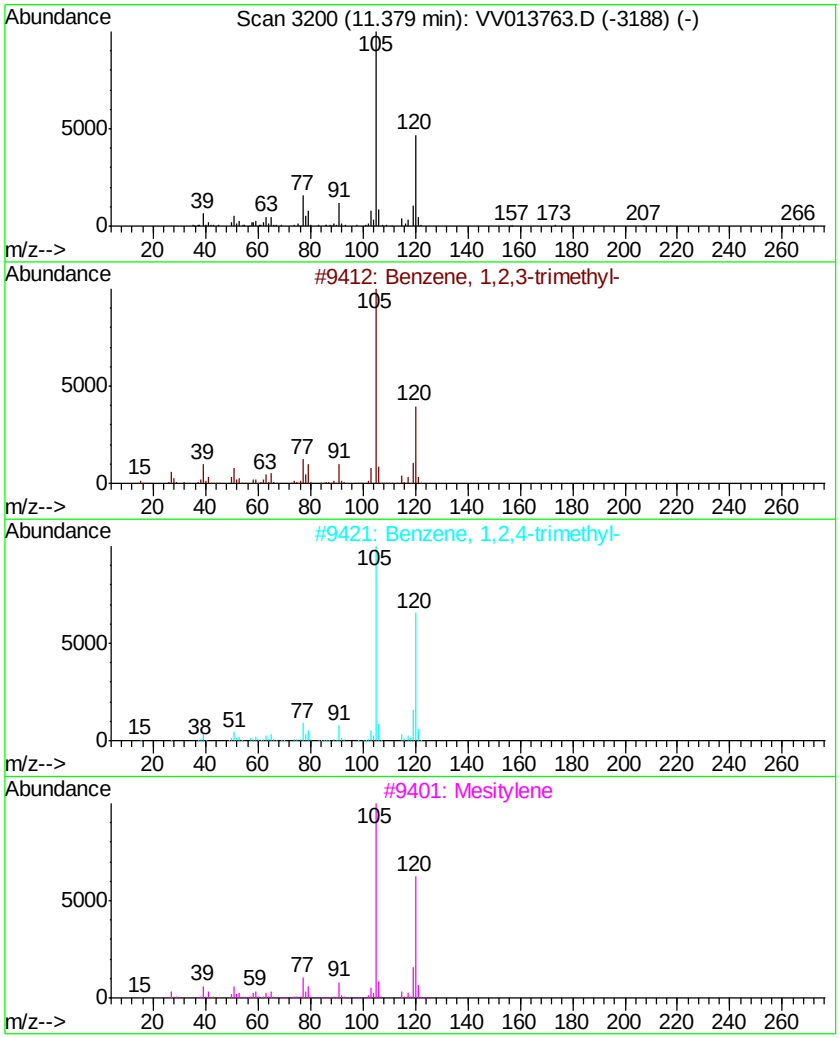
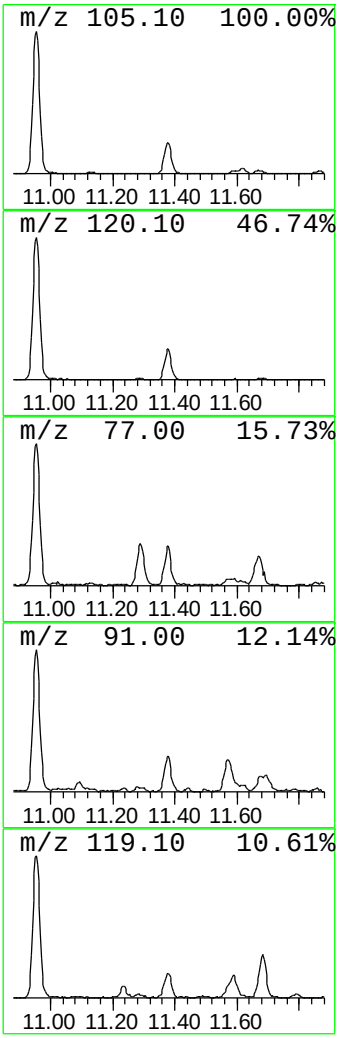
Peak Number 14 Benzene, 1,2,3-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.38	0.65 ug/L	142148	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
2	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
3	Mesitylene	120	C9H12	000108-67-8	94
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\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

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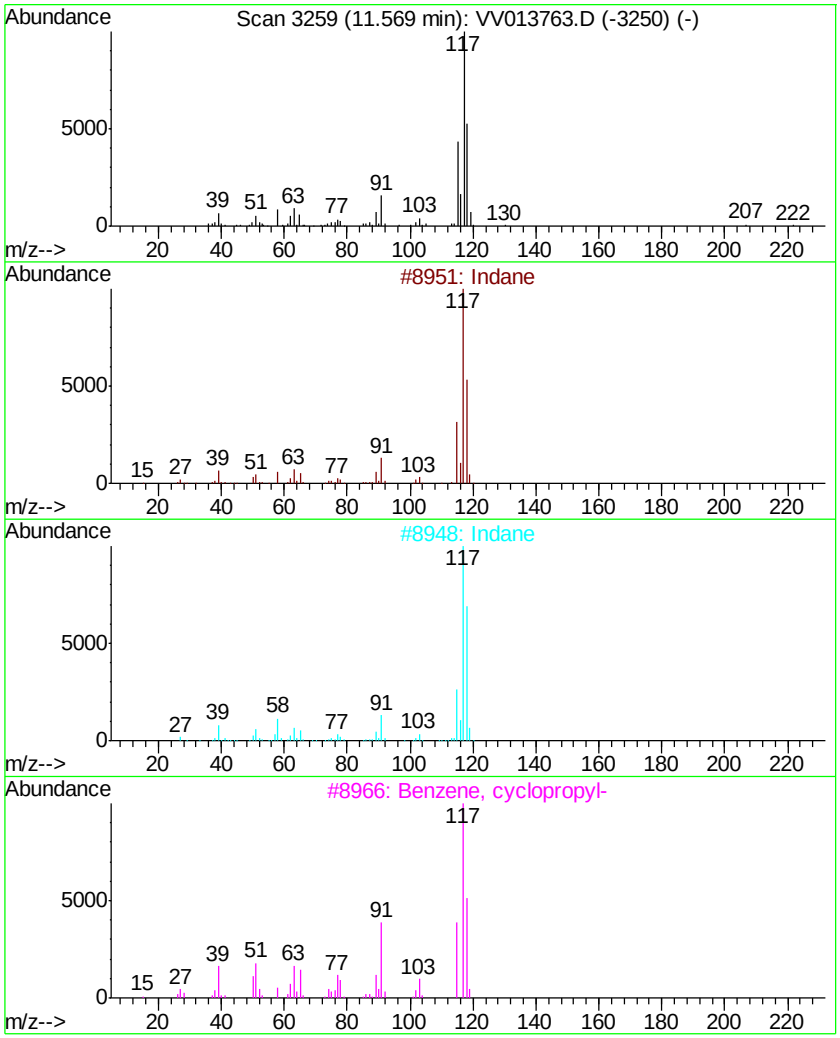
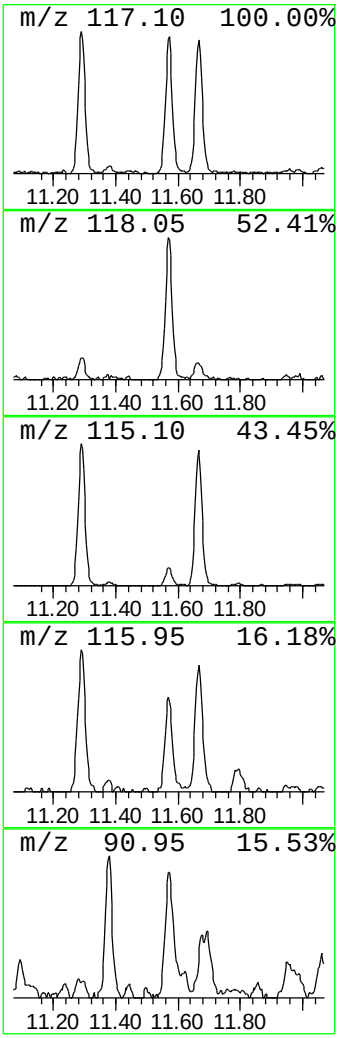
Peak Number 15 Indane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	0.54 ug/L	118209	1,4-Dichlorobenzene-d4	11.29

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	93
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
4	Benzene, (3-chloro-1-propenyl)-	152	C9H9Cl	002687-12-9	59
5	3-Bromo-1-phenyl-1-propene	196	C9H9Br	004392-24-9	59





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_V\DATA\VV112519\
Data File : VV013763.D
Acq On : 25 Nov 2019 14:12
Operator : SY/MD
Sample : K5910-21DL2 400X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_V
ClientSampleId : GAQK4DL2

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR112219WMA.M
Quant Title : TRACE VOA SOM01.0

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Sulfur dioxide	1.21	7.6	ug/L	1274590	1	5.66	842515	5.0
(DEL) Alkane: Str...	1.65	4.3	ug/L	721932	1	5.66	842515	5.0
(DEL) Alkane: Str...	1.82	1.2	ug/L	207879	1	5.66	842515	5.0
unknown-01	2.53	0.6	ug/L	94587	1	5.66	842515	5.0
(DEL) Alkane: Cyc...	2.60	0.6	ug/L	100699	1	5.66	842515	5.0
(DEL) Alkane: Str...	2.79	0.7	ug/L	110285	1	5.66	842515	5.0
(DEL) Alkane: Cyc...	3.80	1.2	ug/L	201286	1	5.66	842515	5.0
Benzene, propyl-	10.39	0.5	ug/L	110116	3	11.29	1097020	5.0
Benzene, 1-ethyl-...	10.49	2.0	ug/L	440044	3	11.29	1097020	5.0
Benzene, 1-ethyl-...	10.78	0.7	ug/L	150376	3	11.29	1097020	5.0
Benzene, 1,2,3-tr...	11.38	0.7	ug/L	142148	3	11.29	1097020	5.0
Indane	11.57	0.5	ug/L	118209	3	11.29	1097020	5.0