LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_R\METHODS\SOMRTR101518WMA.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	3.048	218	225	242	rVB2	322594	976484	1.55%	0.489%
2	3.601	306	316	330	rBV	261229	1007884	1.60%	0.504%
3	4.459	436	457	460	rBV4	189051	813448	1.29%	0.407%
4	6.473	777	788	799	rBV3	429167	1356604	2.15%	0.679%
5	7.197	898	907	923	rBV	356565	952226	1.51%	0.476%
6	7.519	948	960	971	rBV	900456	2749933	4.37%	1.376%
7	7.854	1002	1015	1019	rBV	750929	1792677	2.85%	0.897%
8	7.903	1019	1023	1035	rVB2	875799	1893785	3.01%	0.948%
9	8.456	1107	1114	1124	rBV	883680	1668406	2.65%	0.835%
10	8.894	1179	1186	1191	rBV	471454	989504	1.57%	0.495%
11	8.955	1191	1196	1205	rVB	1272102	2453576	3.90%	1.228%
12	9.971	1355	1363	1370	rBV	1239255	2042260	3.24%	1.022%
13	10.592	1456	1465	1478	rBV	682744	1079608	1.71%	0.540%
14	11.315	1574	1584	1590	rBV	10565971	17070269	27.11%	8.542%
15	11.395	1590	1597	1608	rVV	10168283	14252475	22.63%	7.132%
16	11.498	1608	1614	1632	rVB	11382371	18953107	30.10%	9.484%
17	11.826	1656	1668	1676	rBV	2422786	3567008	5.66%	1.785%
18	12.131	1712	1718	1726	rVB	860114	1169160	1.86%	0.585%
19	12.471	1767	1774	1779	rBV	713983	958790	1.52%	0.480%
20	12.538	1779	1785	1788	rVV	2467096	3571119	5.67%	1.787%
21	12.569	1788	1790	1794	rVV	1511118	1825718	2.90%	0.914%
22	12.617	1794	1798	1810	rVB	2126918	2987695	4.74%	1.495%
23	12.769	1816	1823	1831	rBV	531785	852234	1.35%	0.426%
24	12.922	1842	1848	1853	rBV	6164633	8222126	13.06%	4.114%
25	13.226	1892	1898	1899	rBV	1261886	1727039	2.74%	0.864%
26	13.250	1899	1902	1907	rVB	2561493	3817222	6.06%	1.910%
27	13.420	1920	1930	1941	rBV	35825836	62974319	100.00%	31.511%
28	13.524	1941	1947	1955	rVV2	1708425	3487999	5.54%	1.745%
29	13.871	2000	2004	2009	rBV	708563	1041419	1.65%	0.521%
30	14.150	2043	2050	2058	rVB2	781278	1801964	2.86%	0.902%
31	14.351	2078	2083	2089	rBV	1056657	1418799	2.25%	0.710%
32	14.473	2095	2103	2108	rBV3	1035167	1791942	2.85%	0.897%
33	14.619	2121	2127	2136	rVB	791943	1181364	1.88%	0.591%
34	15.002	2184	2190	2200	rVB	11568170	16343148	25.95%	8.178%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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_R\METHODS\SOMRTR101518WMA.M
Title : TRACE VOA SOM01.0

35	16.006	2348	2355	2367	rVB	4649455	7762555	12.33%	3.884%
36	16.194	2378	2386	2394	rVB	1827811	3295878	5.23%	1.649%

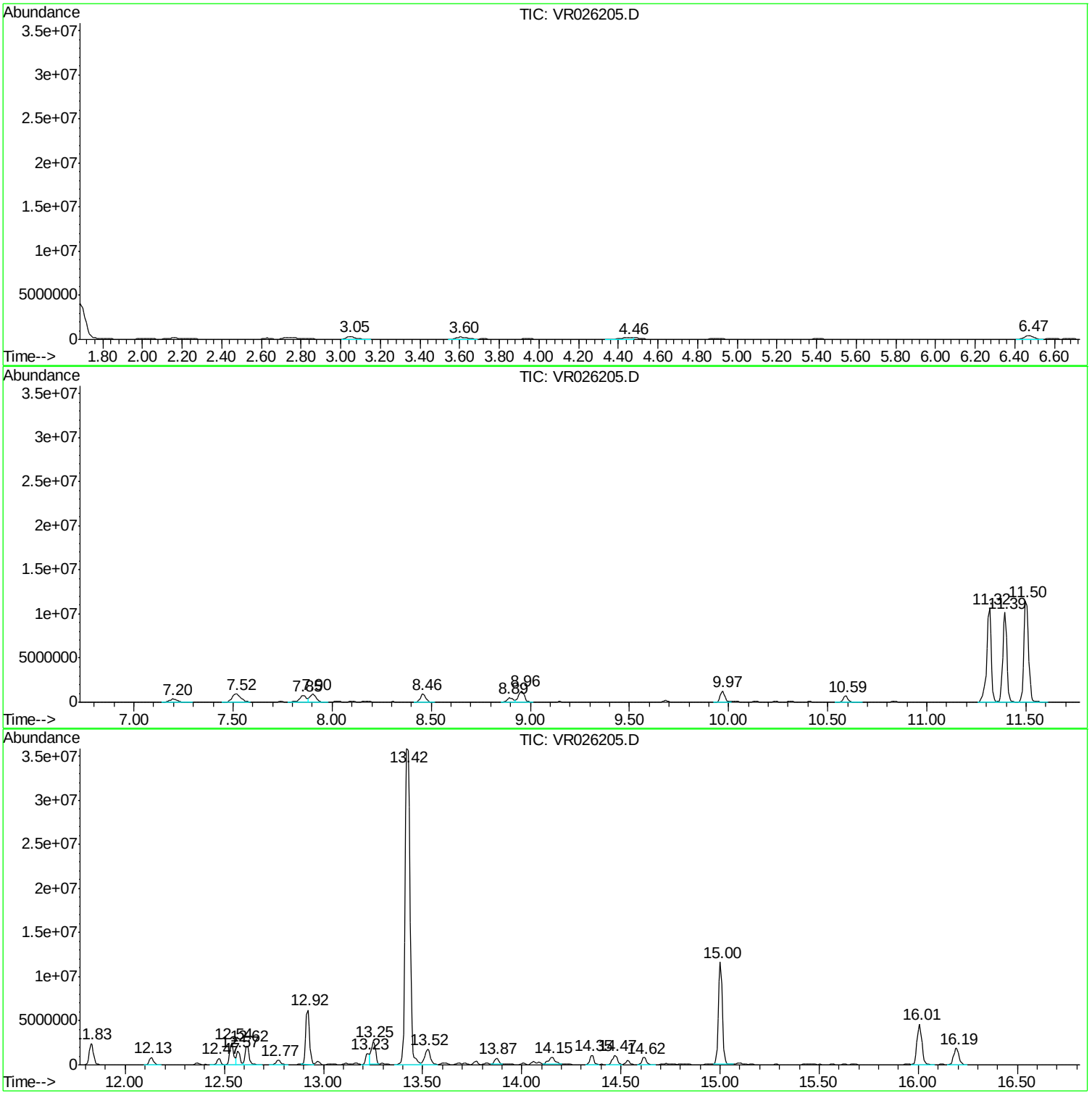
Sum of corrected areas: 199849744

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument :
MSVOA_R
ClientSampleId :
F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

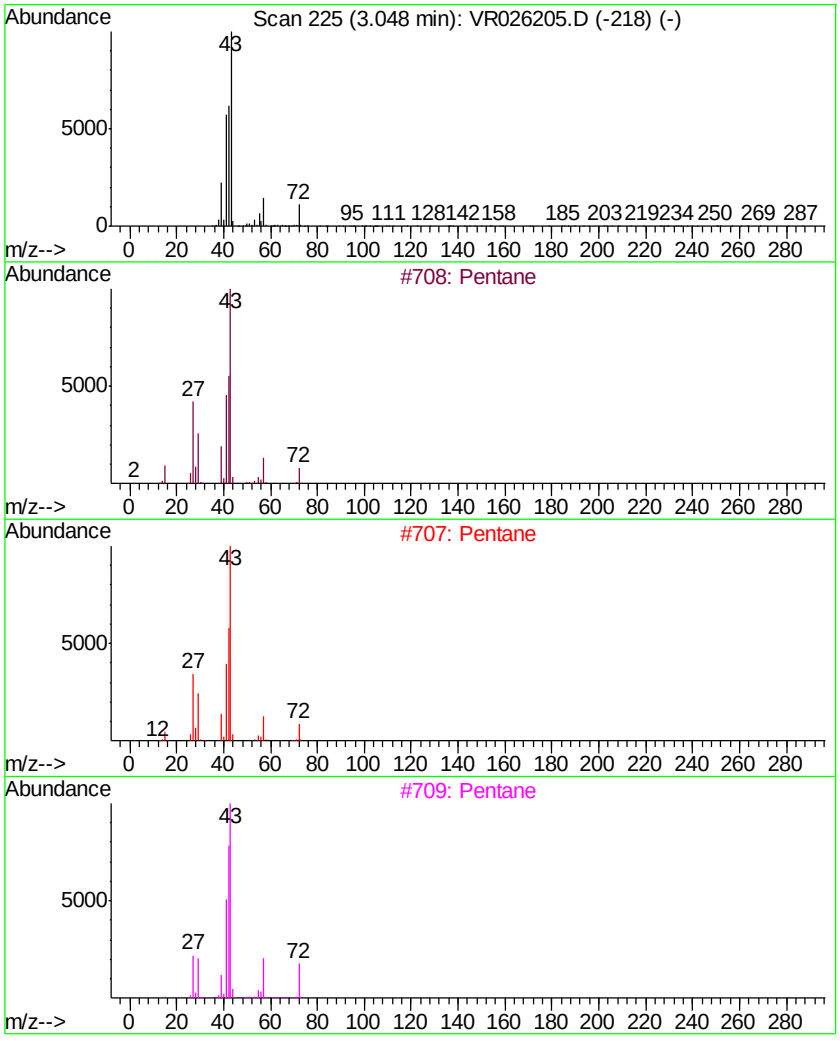
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

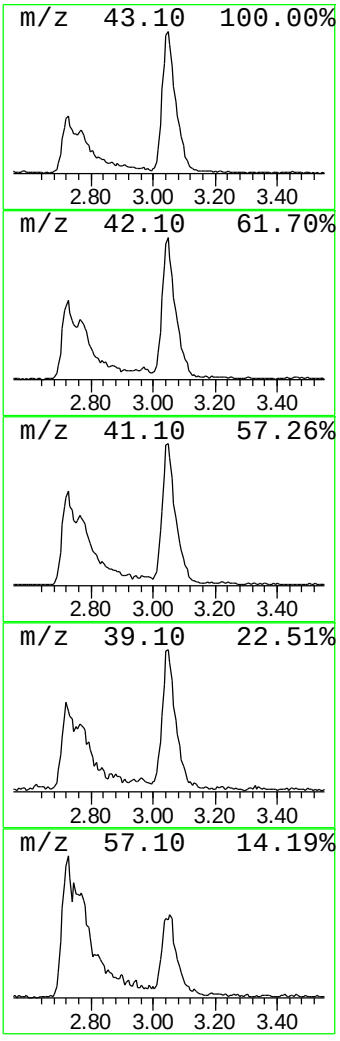
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.05	2.93 ug/L	976484	1,4-Difluorobenzene	8.46

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	59
4	Pentane	72	C5H12	000109-66-0	50
5	Oxirane, ethyl-	72	C4H8O	000106-88-7	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

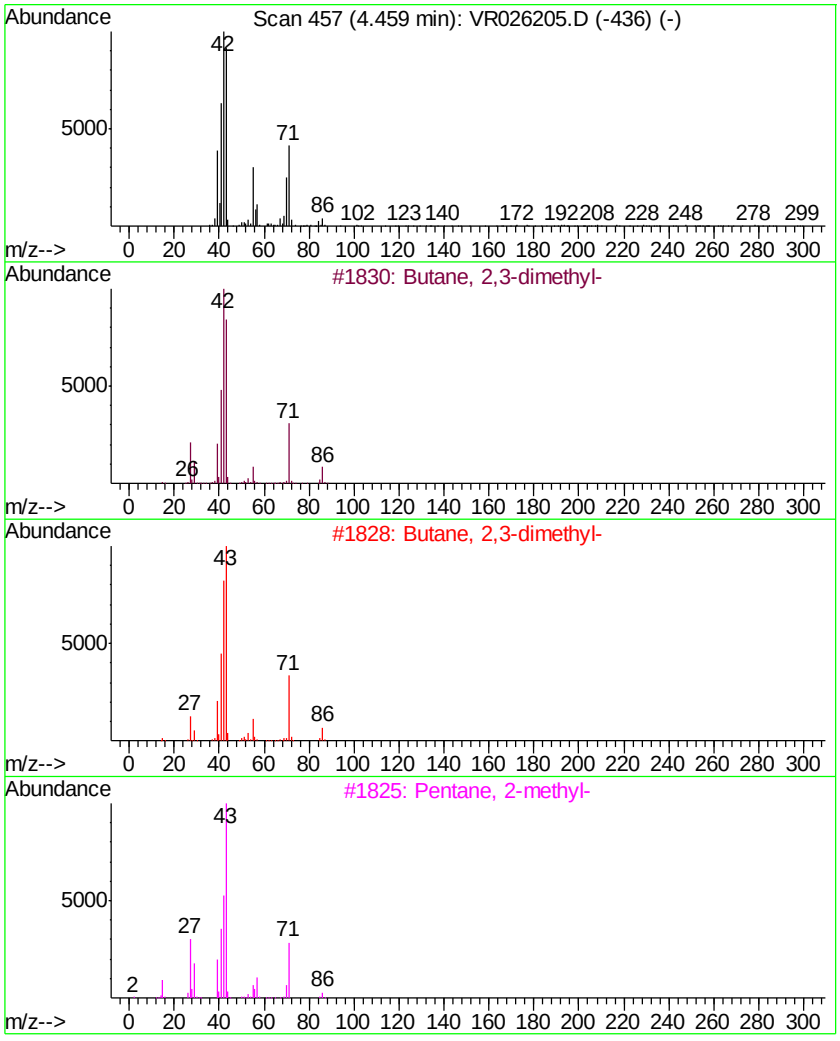
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

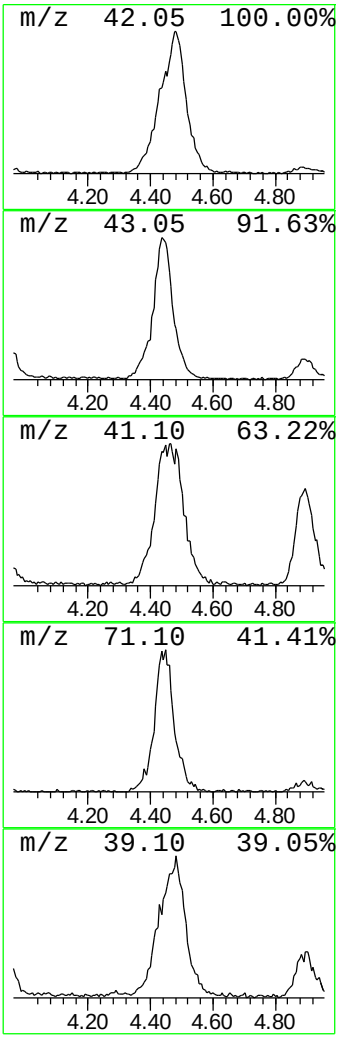
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.46	2.44 ug/L	813448	1,4-Difluorobenzene	8.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	52
2	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	50
3	Pentane, 2-methyl-	86	C6H14	000107-83-5	43
4	Pentane, 2-methyl-	86	C6H14	000107-83-5	38
5	Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	37





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

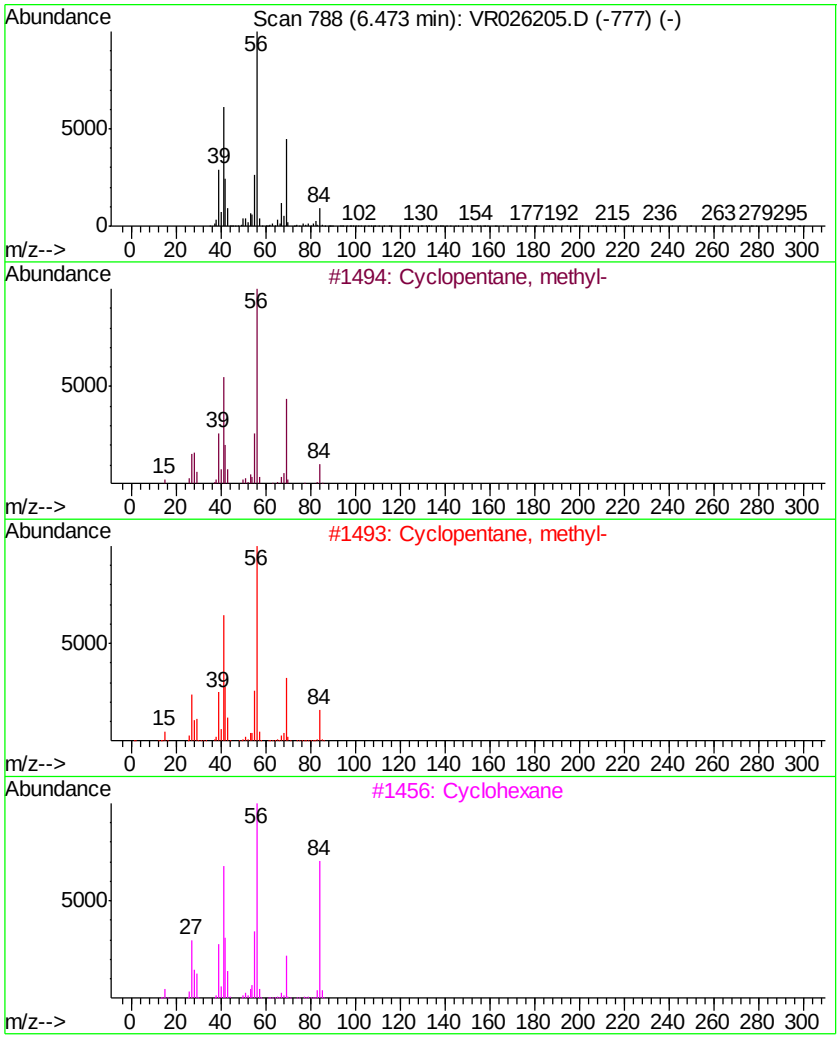
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

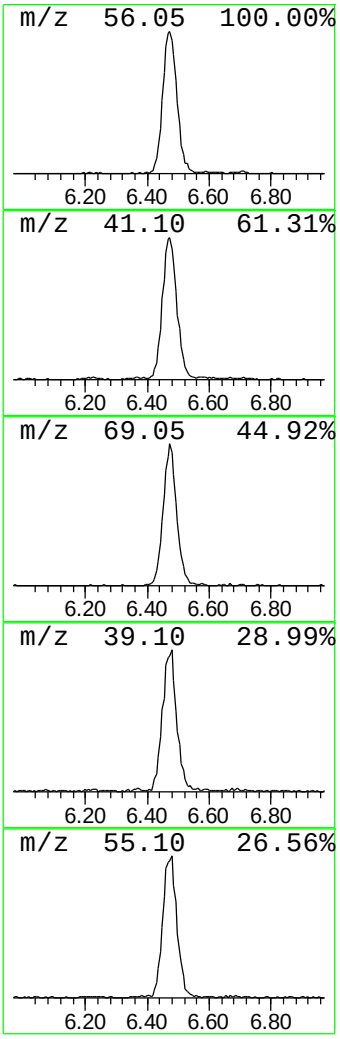
****

Peak Number 3 (DEL) Alkane: Cyclic6.47 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	4.07 ug/L	1356600	1,4-Difluorobenzene	8.46

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	80
3	Cyclohexane	84	C6H12	000110-82-7	72
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
5	Propane, 2-cyclopropyl-	84	C6H12	003638-35-5	59





Library Search Compound Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

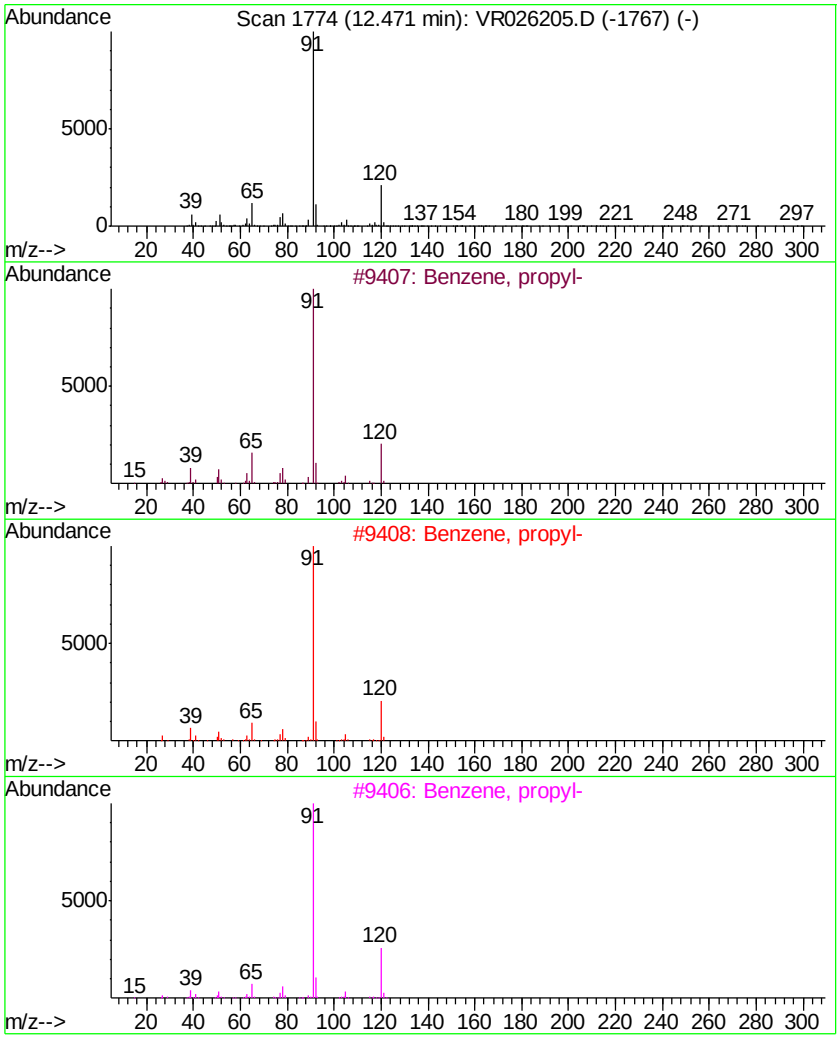
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

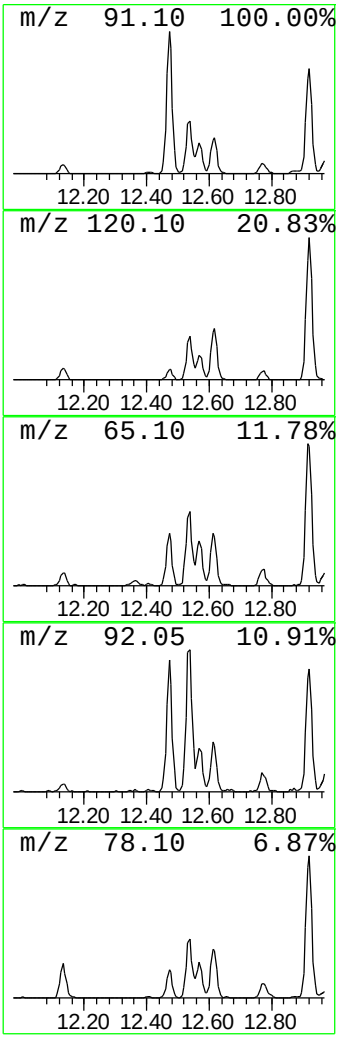
****

Peak Number 4 Benzene, propyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.47	2.78 ug/L	958790	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	95
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	86
4	1,3,5-Cycloheptatriene, 7-ethyl-	120	C9H12	017634-51-4	50
5	Benzeneacetaldehyde	120	C8H8O	000122-78-1	46





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

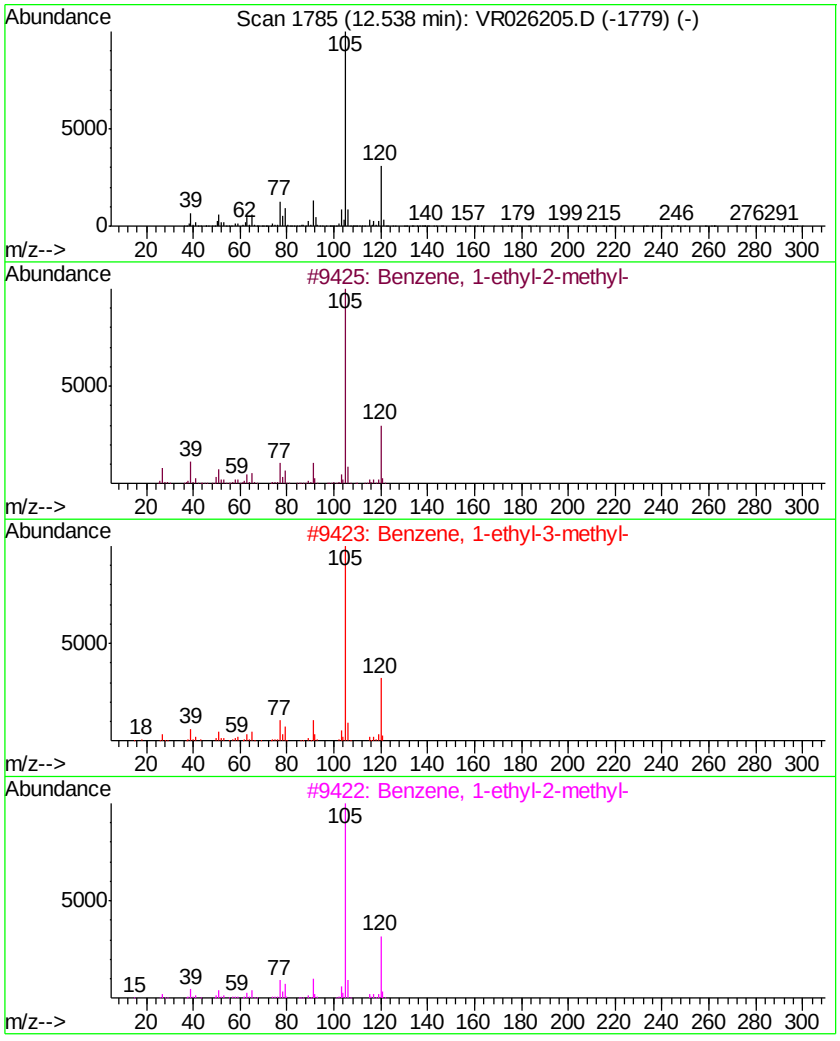
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

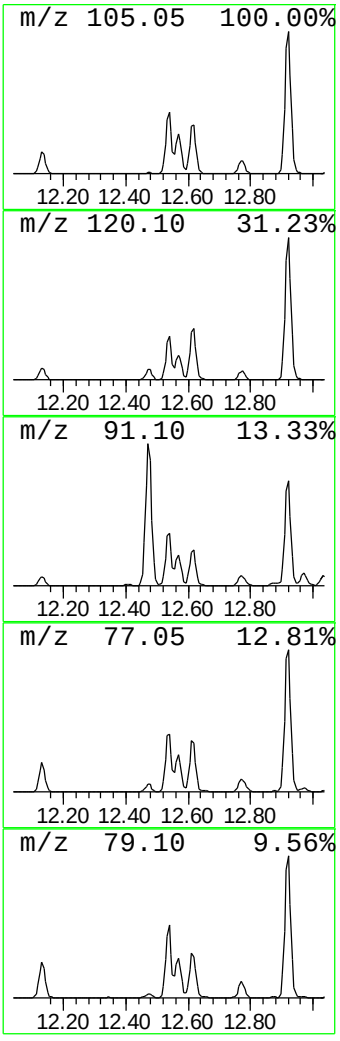
****

Peak Number 5 Benzene, 1-ethyl-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.54	10.34 ug/L	3571120	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	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_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

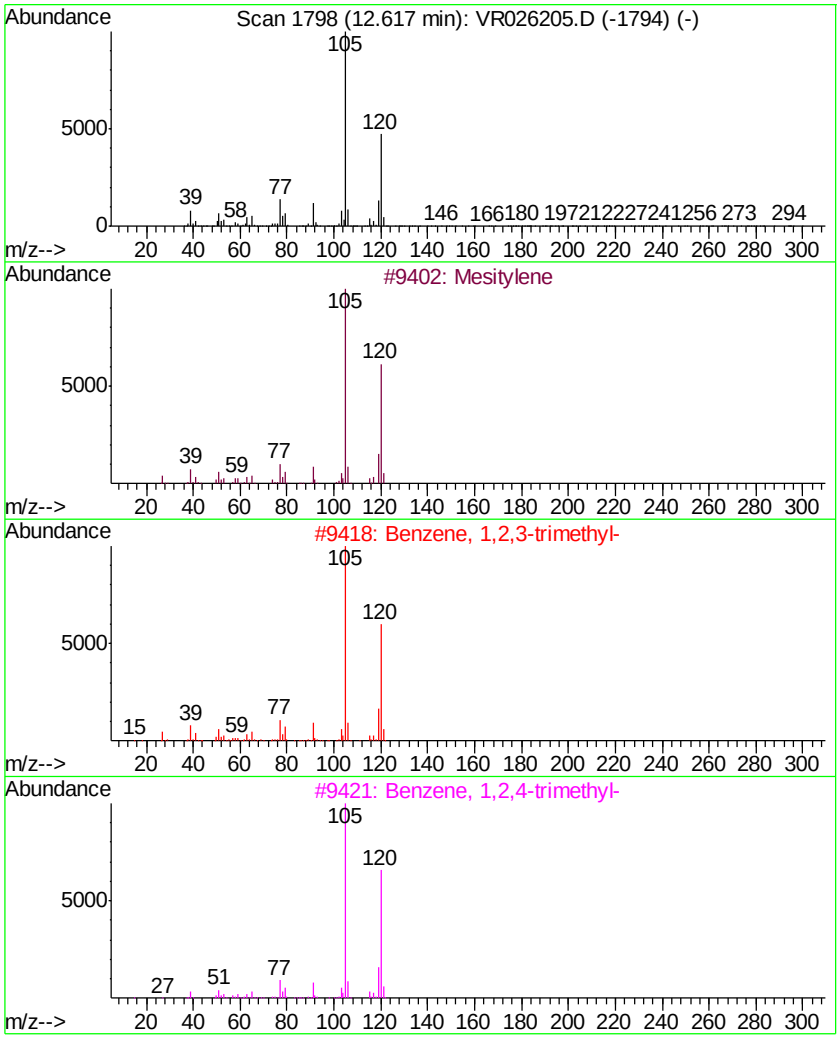
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

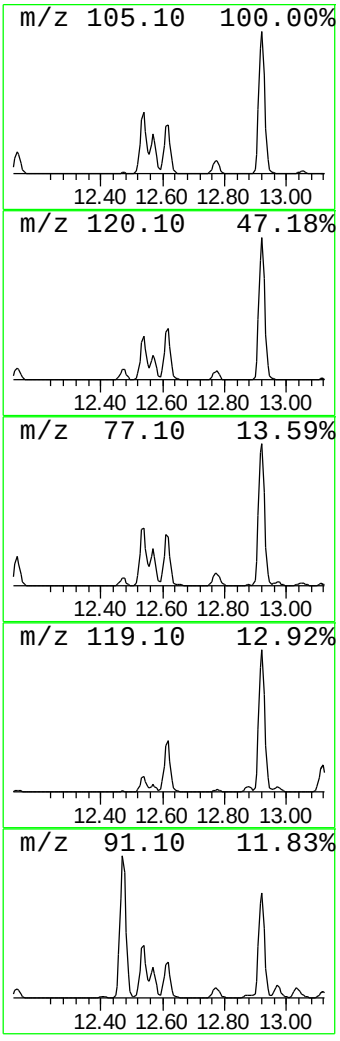
****

Peak Number 7 Mesitylene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.62	8.65 ug/L	2987700	1,4-Dichlorobenzene-d4	13.23

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_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

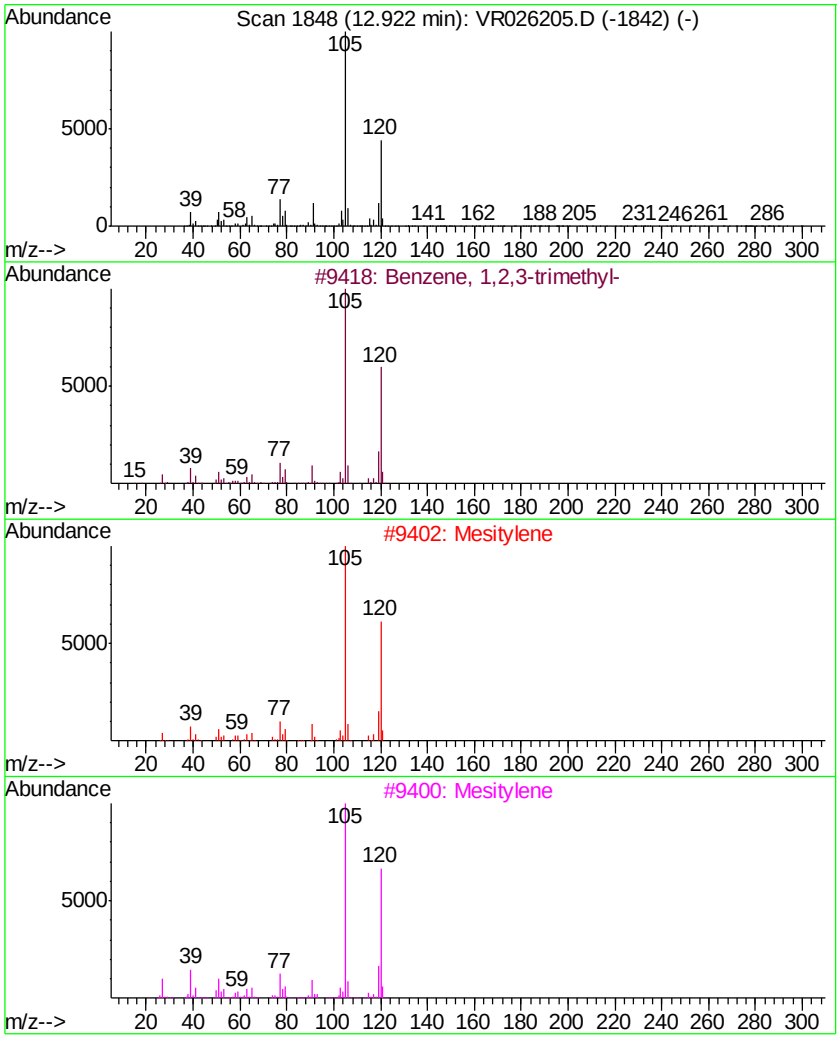
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

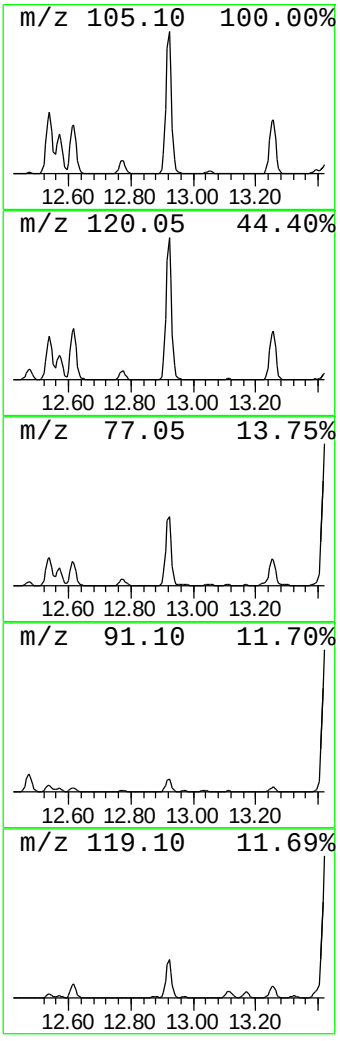
****

Peak Number 9 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.92	23.80 ug/L	8222130	1,4-Dichlorobenzene-d4	13.23

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

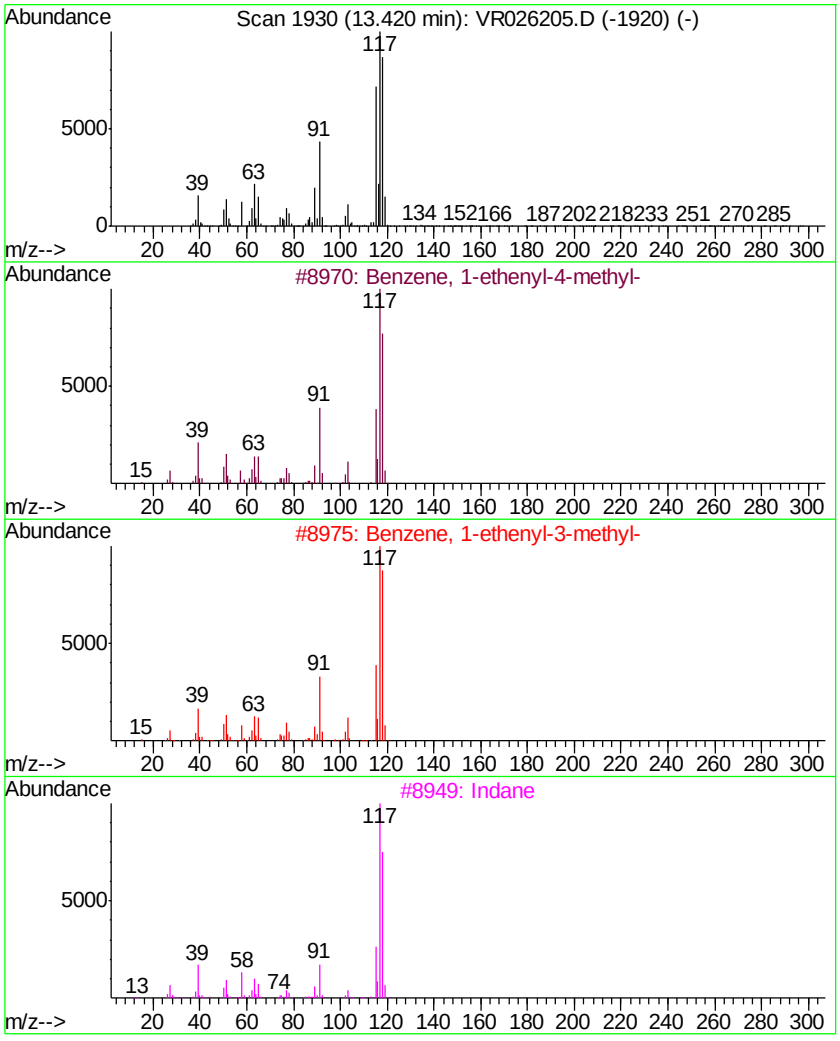
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

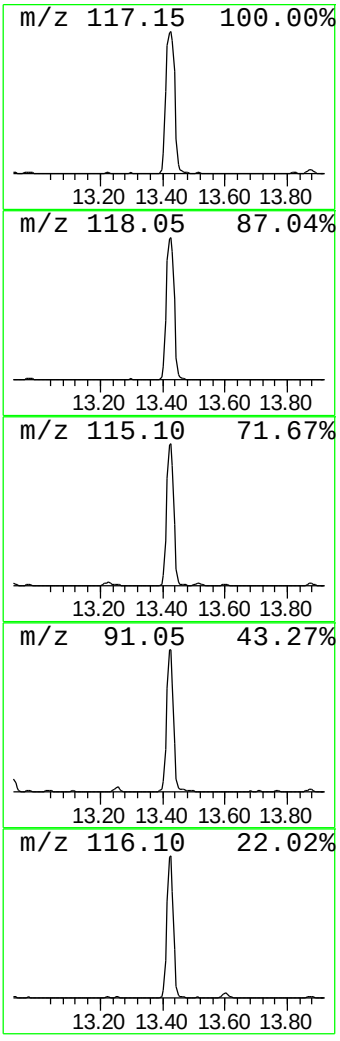
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 1-ethenyl-4-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.42	182.32 ug/L	62974300	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	92
2	Benzene, 1-ethenyl-3-methyl-	118	C9H10	000100-80-1	92
3	Indane	118	C9H10	000496-11-7	81
4	Indane	118	C9H10	000496-11-7	76
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	70





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

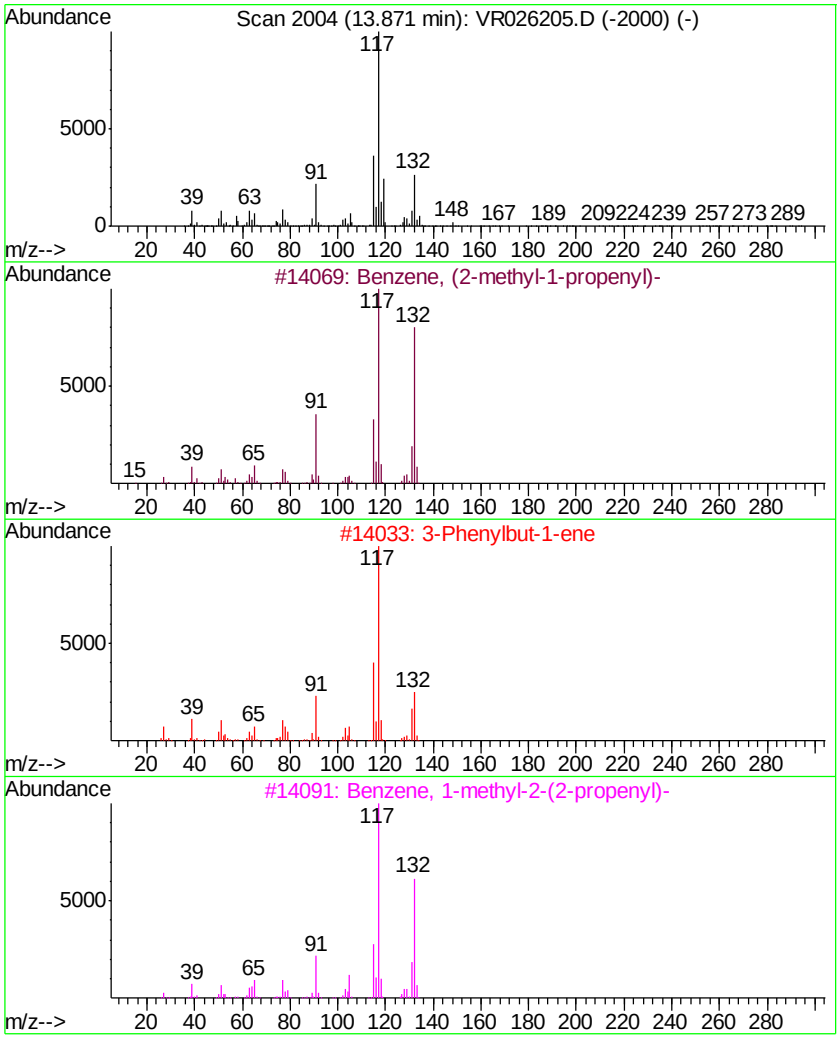
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

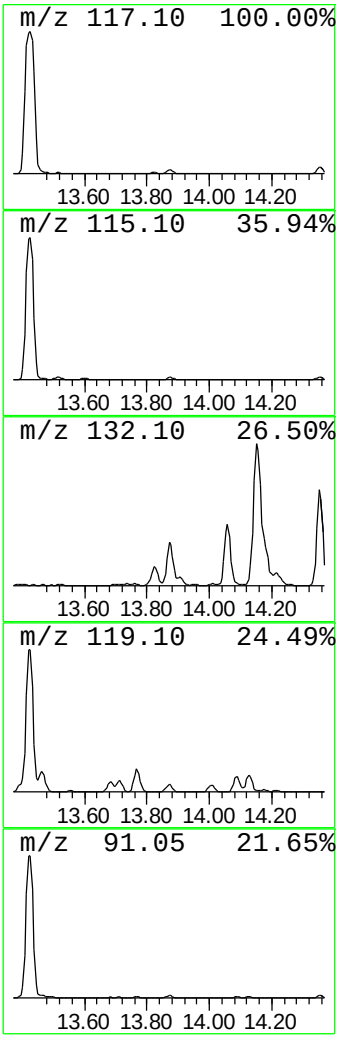
****

Peak Number 11 Benzene, (2-methyl-1-propen... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.87	3.02 ug/L	1041420	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	87
2	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
4	Indan, 1-methyl-	132	C10H12	000767-58-8	76
5	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	74





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

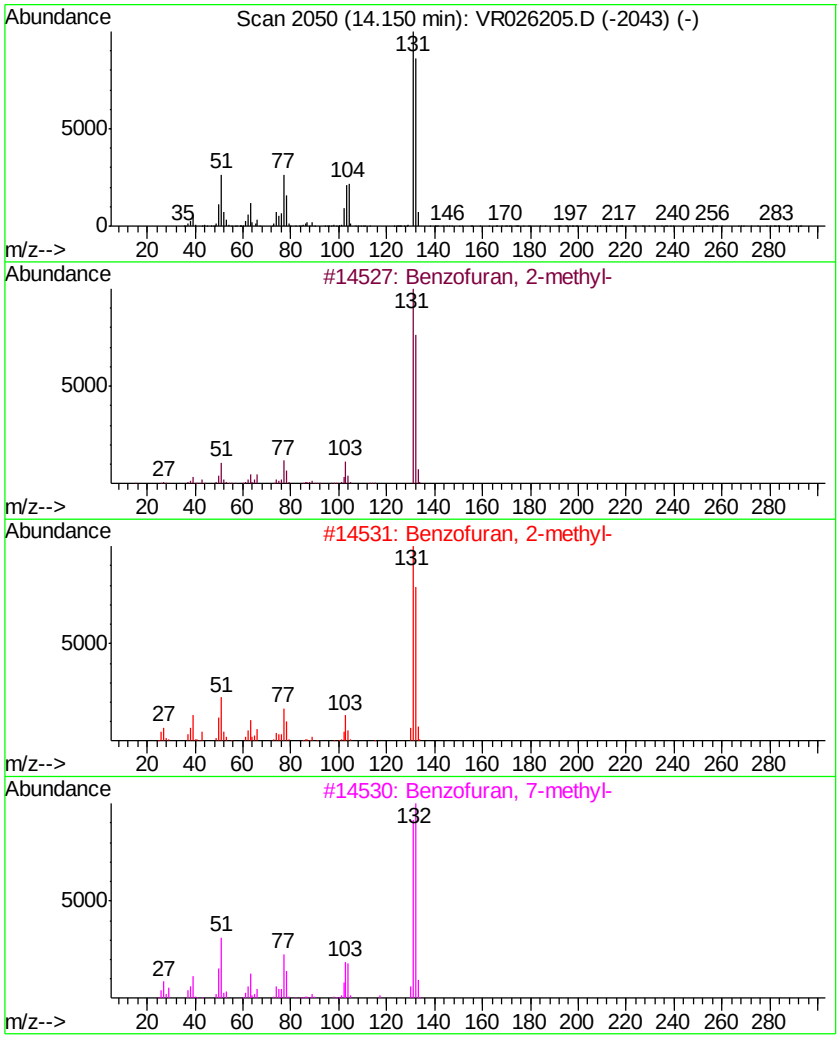
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

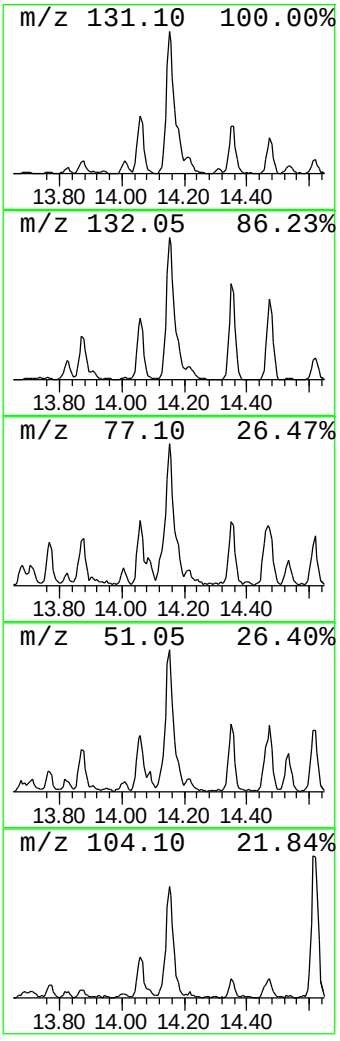
****

Peak Number 12 Benzofuran, 2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.15	5.22 ug/L	1801960	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	95
2	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3	Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	93
4	Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
5	3-Phenyl-2-propyn-1-ol	132	C9H8O	001504-58-1	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

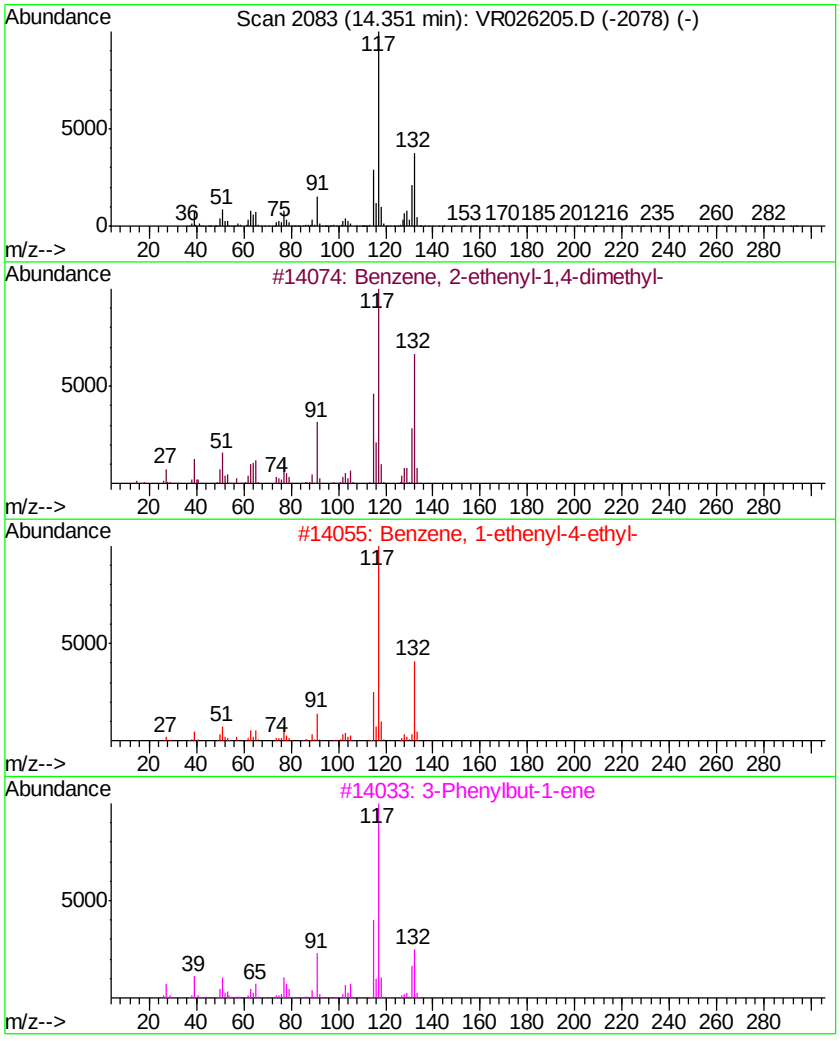
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

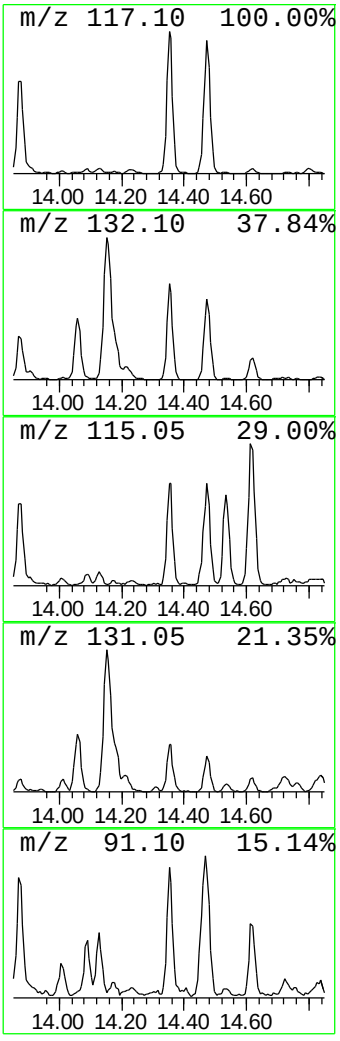
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.35	4.11 ug/L	1418800	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	91
4	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

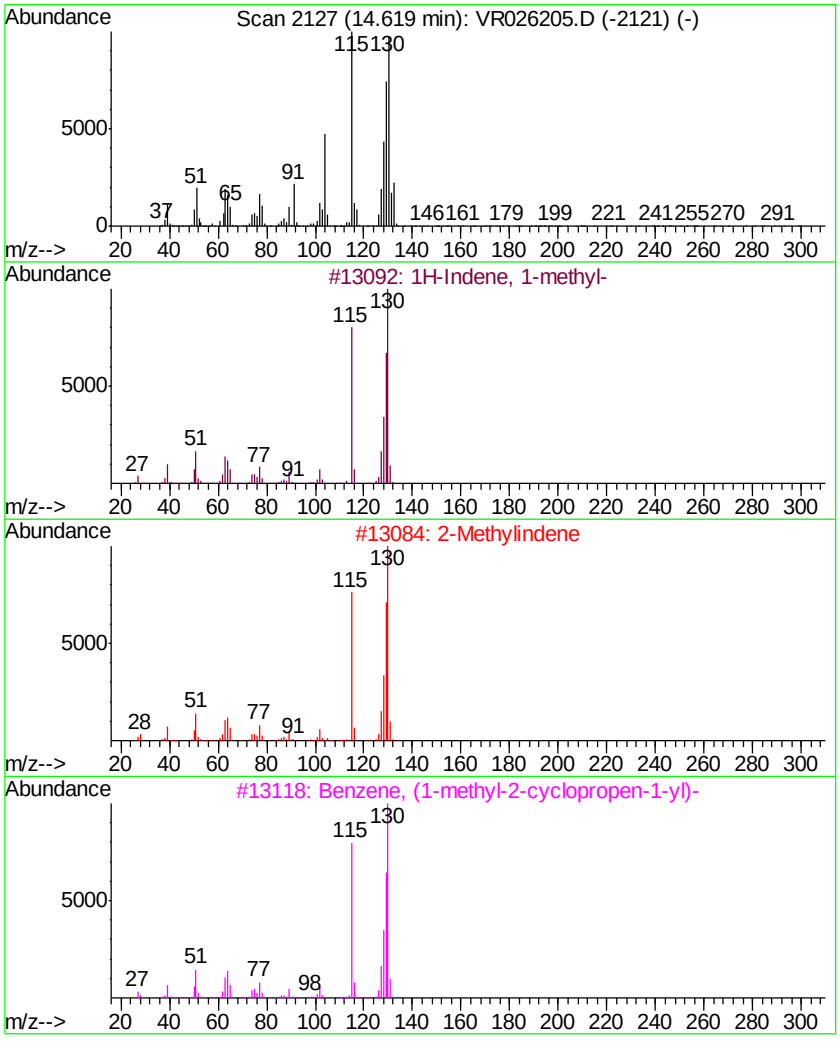
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

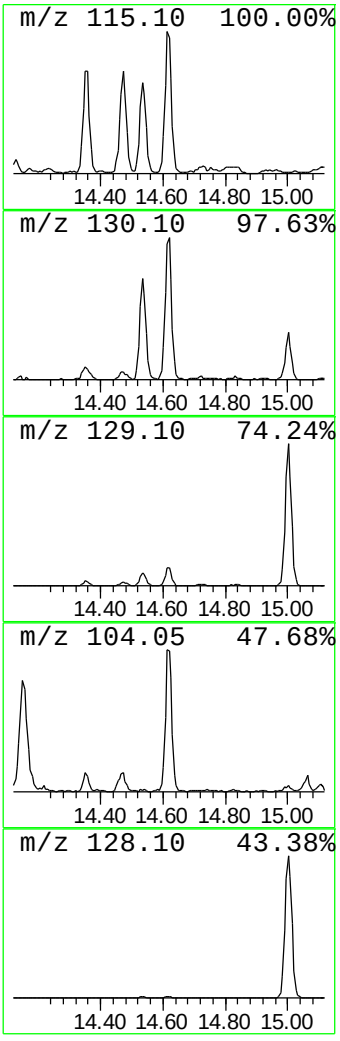
****

Peak Number 15 1H-Indene, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.62	3.42 ug/L	1181360	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	96
2	2-Methylindene	130	C10H10	002177-47-1	95
3	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	93
4	1H-Indene, 3-methyl-	130	C10H10	000767-60-2	91
5	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.M
Quant Title : TRACE VOA SOM01.0

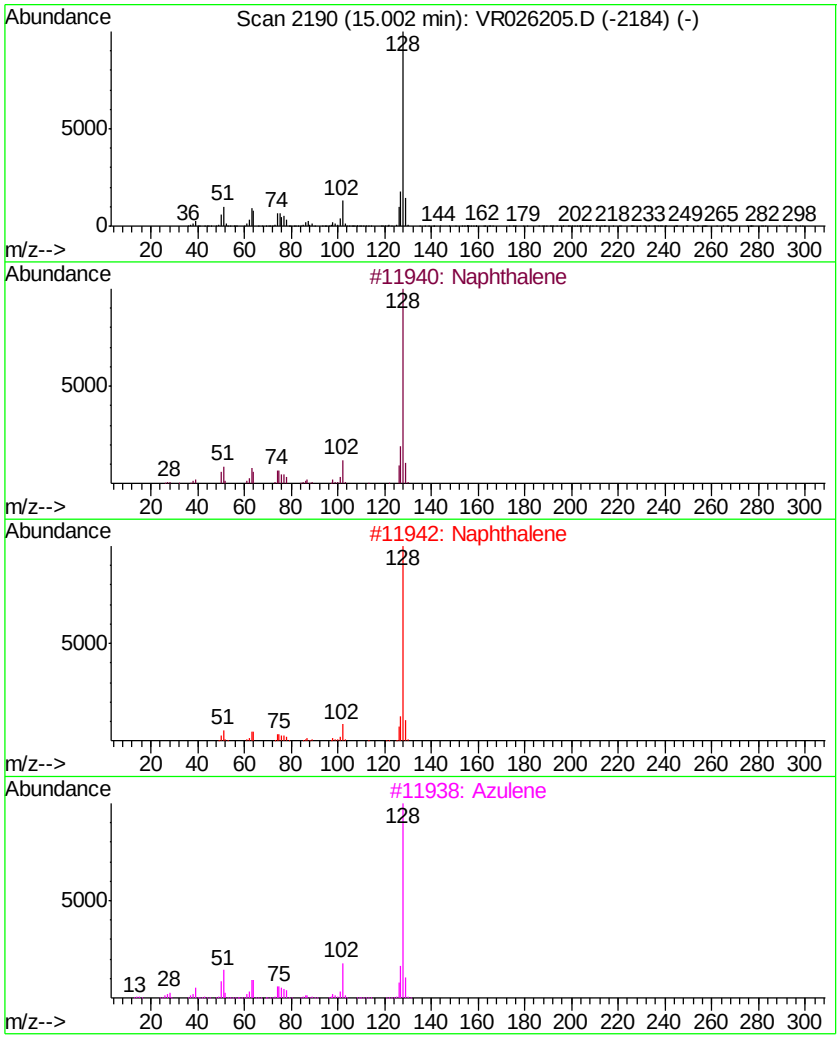
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

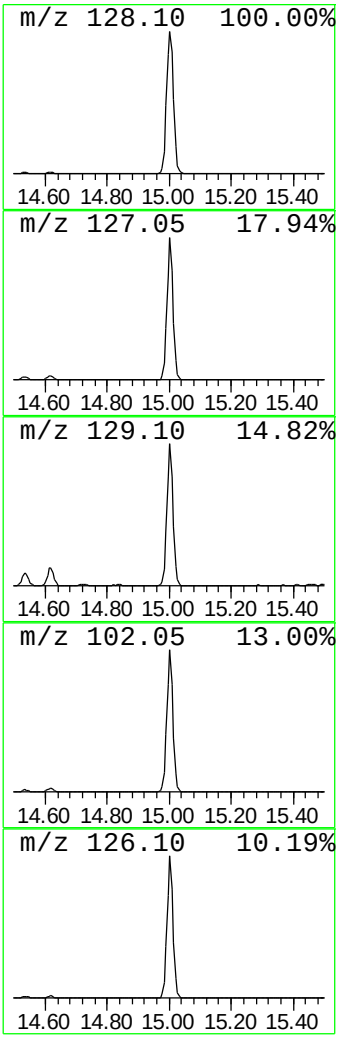
****

Peak Number 16 Azulene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.00	47.32 ug/L	16343100	1,4-Dichlorobenzene-d4	13.23

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	96
2	Naphthalene	128	C10H8	000091-20-3	95
3	Azulene	128	C10H8	000275-51-4	94
4	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	91
5	Naphthalene	128	C10H8	000091-20-3	91





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_R\Data\VR110718\
Data File : VR026205.D
Acq On : 7 Nov 2018 15:36
Operator : SY/MD
Sample : J5801-06
Misc : 25mL/MSVOA_R/WATER
ALS Vial : 10 Sample Multiplier: 1

Instrument : MSVOA_R
ClientSampleId : F9D05

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_R\METHODS\SOMRTR101518WMA.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
(DEL) Alkane: Str...	3.05	2.9	ug/L	976484	1	8.46	1668410	5.0
(DEL) Alkane: Str...	4.46	2.4	ug/L	813448	1	8.46	1668410	5.0
(DEL) Alkane: Cyc...	6.47	4.1	ug/L	1356600	1	8.46	1668410	5.0
Benzene, propyl-	12.47	2.8	ug/L	958790	3	13.23	1727040	5.0
Benzene, 1-ethyl-...	12.54	10.3	ug/L	3571120	3	13.23	1727040	5.0
Mesitylene	12.62	8.7	ug/L	2987700	3	13.23	1727040	5.0
Benzene, 1,2,3-tr...	12.92	23.8	ug/L	8222130	3	13.23	1727040	5.0
Benzene, 1-etheny...	13.42	182.3	ug/L	62974300	3	13.23	1727040	5.0
Benzene, (2-methy...	13.87	3.0	ug/L	1041420	3	13.23	1727040	5.0
Benzofuran, 2-met...	14.15	5.2	ug/L	1801960	3	13.23	1727040	5.0
Benzene, 2-etheny...	14.35	4.1	ug/L	1418800	3	13.23	1727040	5.0
1H-Indene, 1-methyl-	14.62	3.4	ug/L	1181360	3	13.23	1727040	5.0
Azulene	15.00	47.3	ug/L	16343100	3	13.23	1727040	5.0