LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

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_N\methods\82N122721W.M
Title : SW846 8260

Signal : TIC: VN070437.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.292	80	113	133	rBV	1966451	9790724	2.95%	0.827%
2	3.140	413	429	467	rVB	4891434	11463022	3.45%	0.968%
3	3.513	553	568	625	rVB2	2023876	6613302	1.99%	0.558%
4	6.900	1810	1831	1865	rVB	1355800	3973018	1.20%	0.335%
5	7.147	1899	1923	1943	rBV	3801804	11094890	3.34%	0.937%
6	7.841	2165	2182	2203	rVB	1420053	3437948	1.04%	0.290%
7	8.088	2259	2274	2293	rVV3	2560990	6844431	2.06%	0.578%
8	8.965	2579	2601	2630	rBV3	2795369	7010455	2.11%	0.592%
9	10.440	3133	3151	3166	rBV2	3618467	7012337	2.11%	0.592%
10	11.744	3618	3637	3653	rBV	2894076	5409379	1.63%	0.457%
11	11.843	3655	3674	3694	rBV2	51163869	94782229	28.55%	8.002%
12	11.956	3695	3716	3759	rVB6	144226617	331982695	100.00%	28.029%
13	12.277	3812	3836	3855	rBV	12008538	20426240	6.15%	1.725%
14	12.575	3935	3947	3967	rVB	4232888	7097927	2.14%	0.599%
15	12.728	3990	4004	4025	rBV	2630057	4432402	1.34%	0.374%
16	12.918	4060	4075	4085	rBV	9272342	15240514	4.59%	1.287%
17	12.980	4085	4098	4105	rBV2	41539538	70217666	21.15%	5.928%
18	13.058	4119	4127	4147	rVB	23314067	37157250	11.19%	3.137%
19	13.219	4170	4187	4211	rVB2	35074568	59591872	17.95%	5.031%
20	13.366	4216	4242	4271	rBV3	90578194	164862210	49.66%	13.919%
21	13.702	4343	4367	4403	rVB2	43457057	78975056	23.79%	6.668%
22	13.873	4404	4431	4467	rBV4	45378408	93096386	28.04%	7.860%
23	14.128	4514	4526	4532	rBV	3986132	6503457	1.96%	0.549%
24	14.214	4547	4558	4571	rVB	6262238	9753703	2.94%	0.823%
25	14.324	4589	4599	4620	rVB2	5070742	8716056	2.63%	0.736%
26	14.538	4663	4679	4685	rBV	3741120	6016378	1.81%	0.508%
27	14.576	4685	4693	4705	rVB	5150868	8004051	2.41%	0.676%
28	14.807	4767	4779	4791	rBV	4608471	7286391	2.19%	0.615%
29	14.927	4805	4824	4838	rBV3	9096128	18311218	5.52%	1.546%
30	15.501	5008	5038	5062	rBV2	32354115	64048899	19.29%	5.408%
31	16.614	5436	5453	5481	rVB	2352310	5278210	1.59%	0.446%

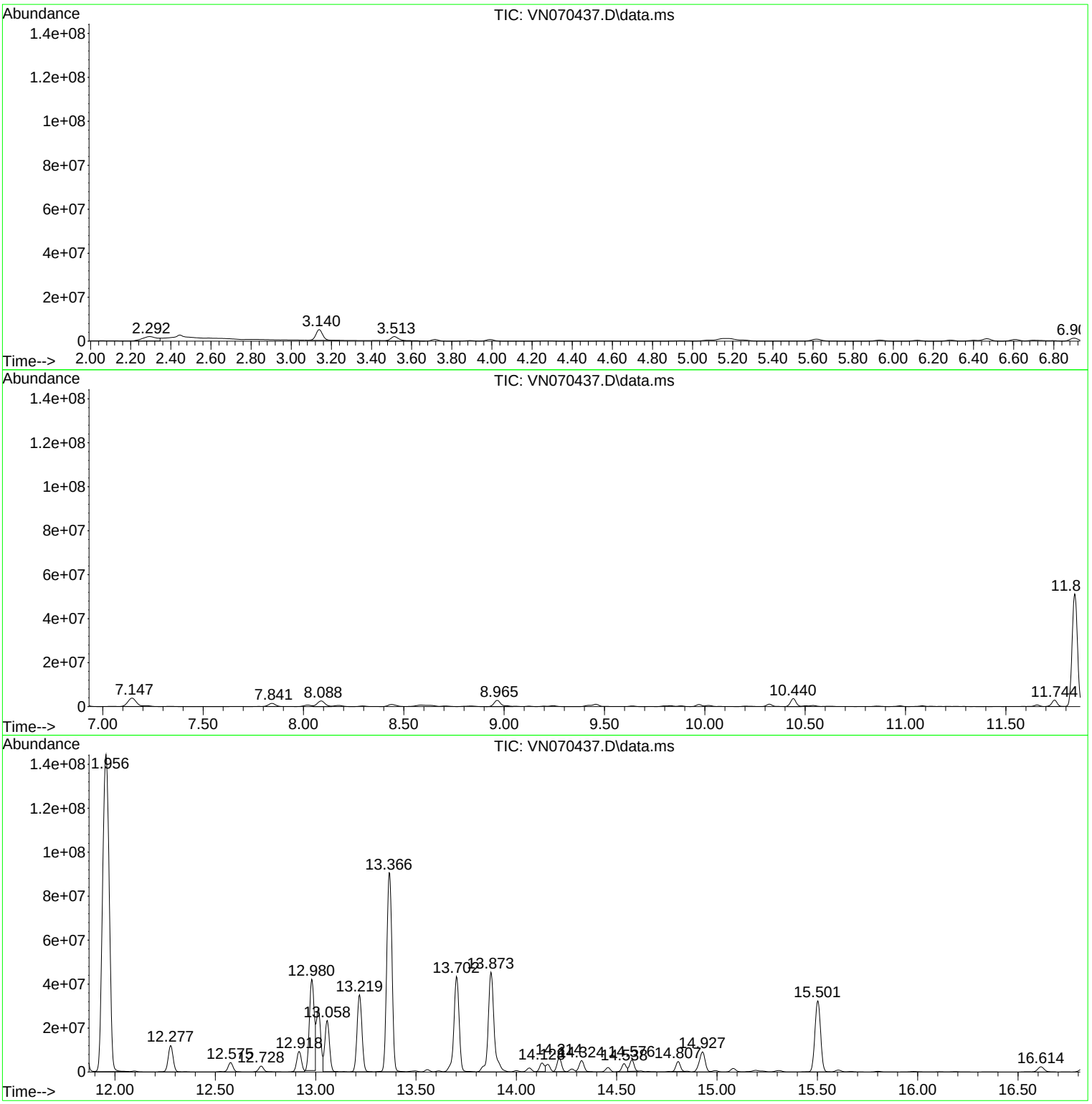
Sum of corrected areas: 1184430316

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

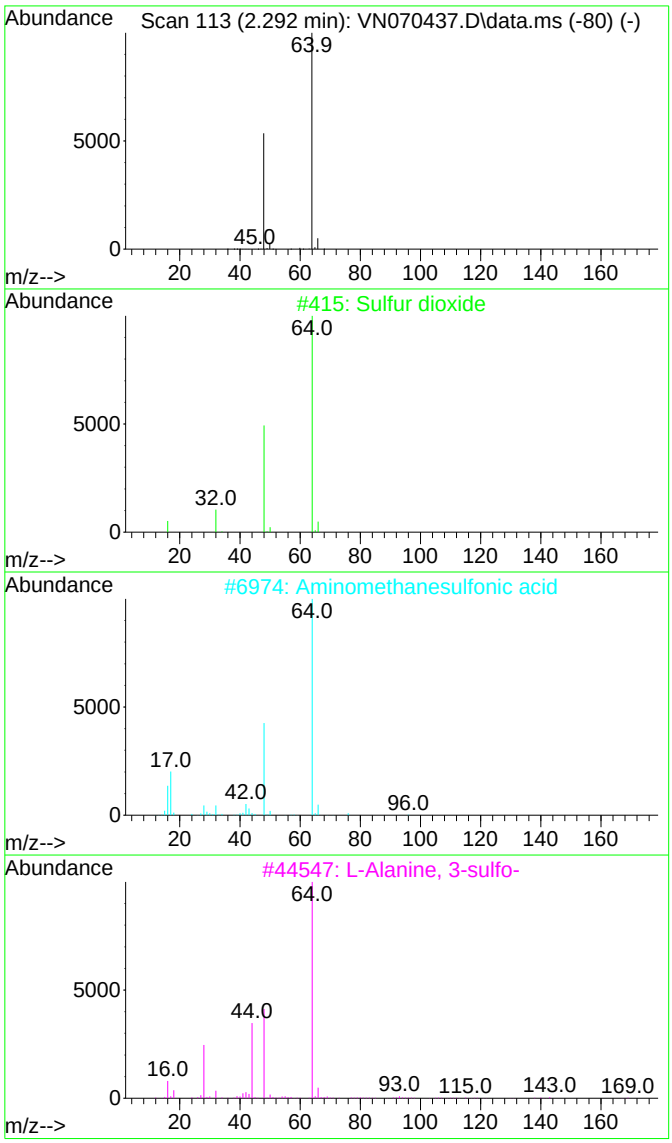
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

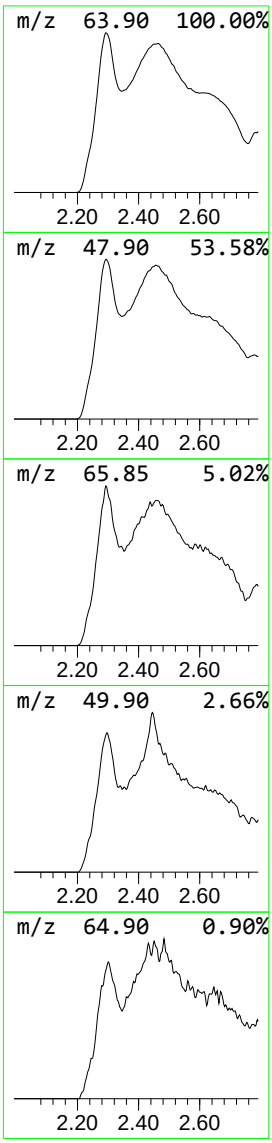
****

Peak Number 1 Sulfur dioxide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.292	71.52 ug/l	9790720	Pentafluorobenzene	8.086

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	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4
5	Ethyl Chloride	64	C2H5Cl	000075-00-3	3





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

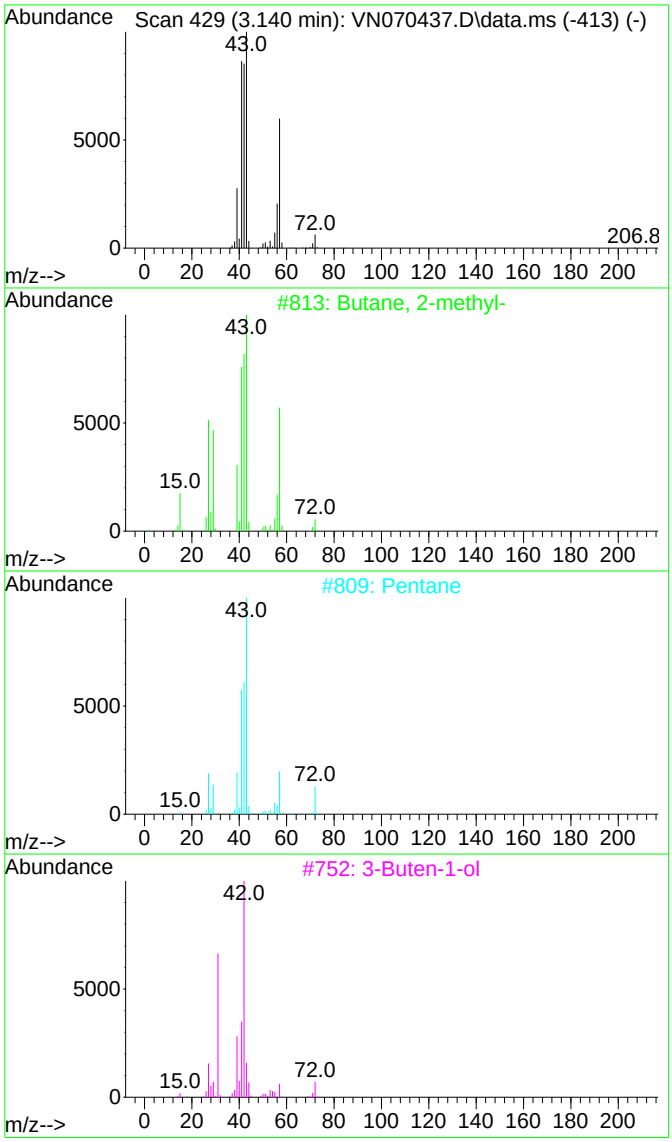
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

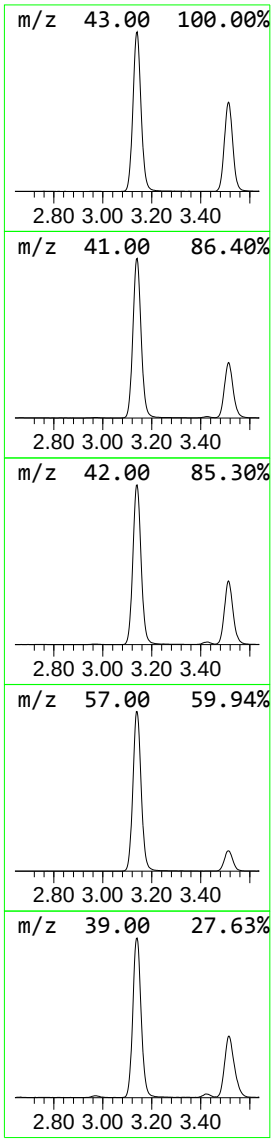
****

Peak Number 2 Butane, 2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.140	83.74 ug/l	11463000	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	90
2	Pentane	72	C5H12	000109-66-0	33
3	3-Buten-1-ol	72	C4H8O	000627-27-0	25
4	1-Propene, 2-methyl-	56	C4H8	000115-11-7	9
5	Methane, isocyanato-	57	C2H3NO	000624-83-9	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

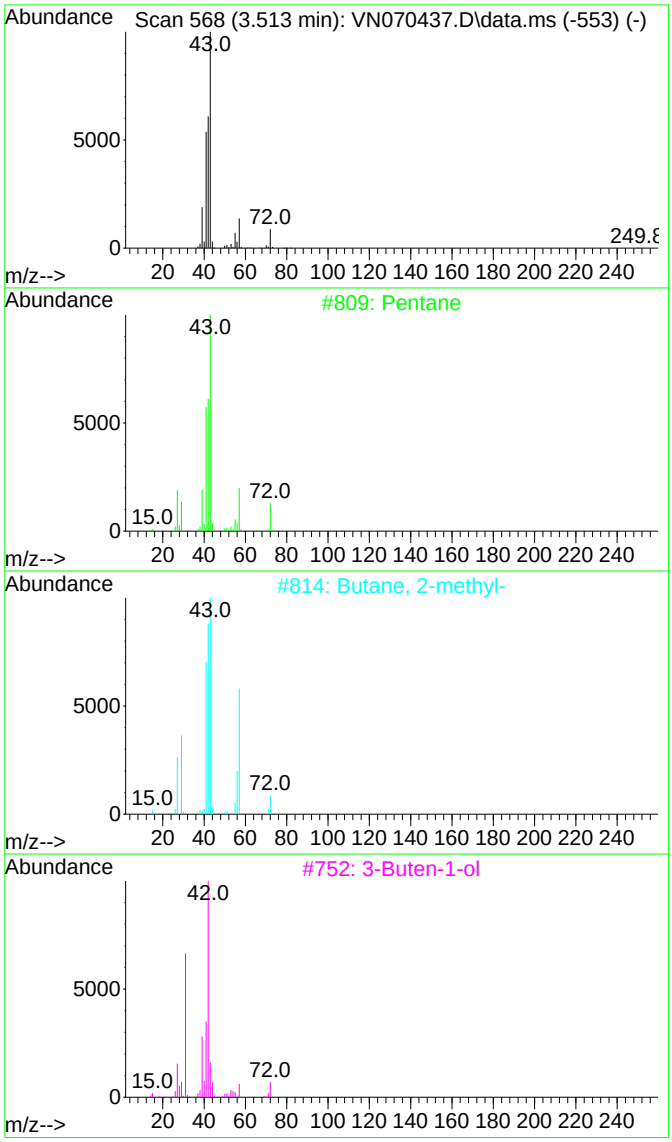
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

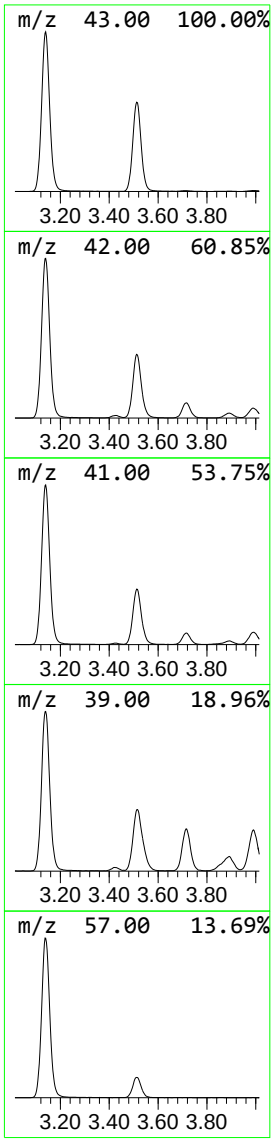
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Pentane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.513	48.31 ug/l	6613300	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Butane, 2-methyl-	72	C5H12	000078-78-4	50
3	3-Buten-1-ol	72	C4H8O	000627-27-0	9
4	Oxirane, ethyl-	72	C4H8O	000106-88-7	9
5	Methyl glyoxal	72	C3H4O2	000078-98-8	9





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

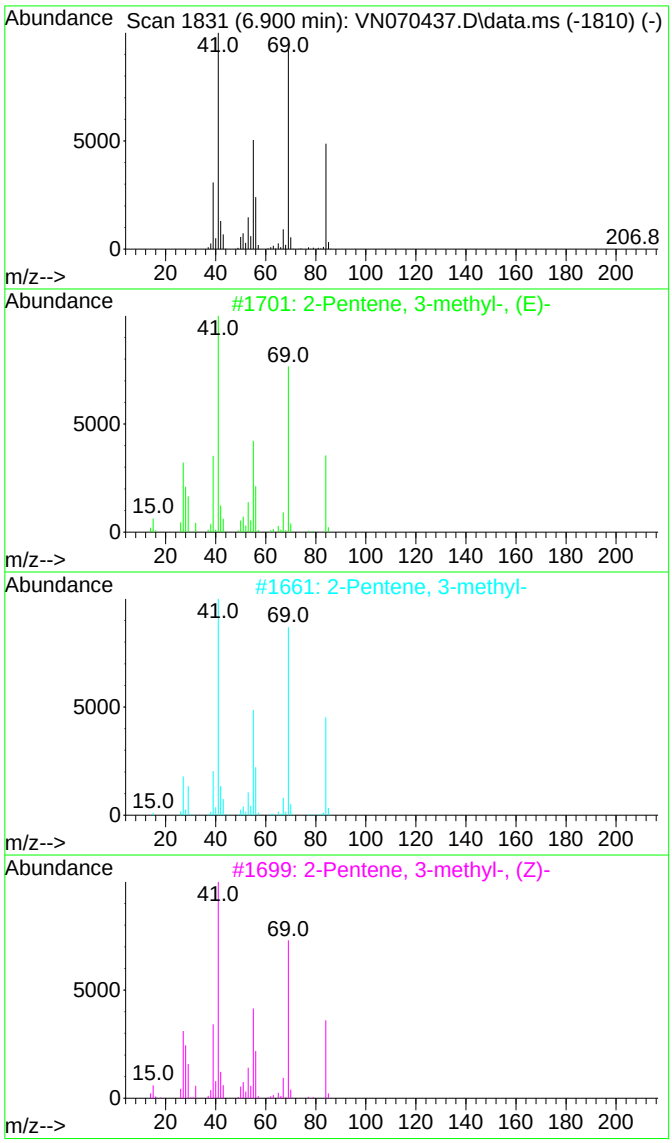
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

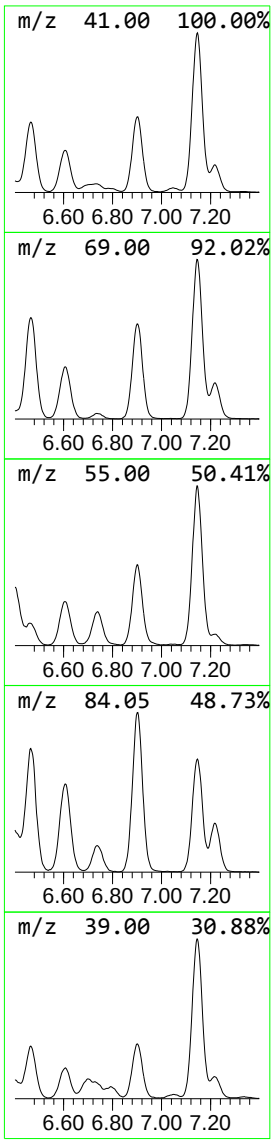
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 2-Pentene, 3-methyl-, (E)- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.900	29.02 ug/l	3973020	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 3-methyl-, (E)-	84	C6H12	000616-12-6	95
2	2-Pentene, 3-methyl-	84	C6H12	000922-61-2	94
3	2-Pentene, 3-methyl-, (Z)-	84	C6H12	000922-62-3	94
4	Cyclopropane, 1,1,2-trimethyl-	84	C6H12	004127-45-1	90
5	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

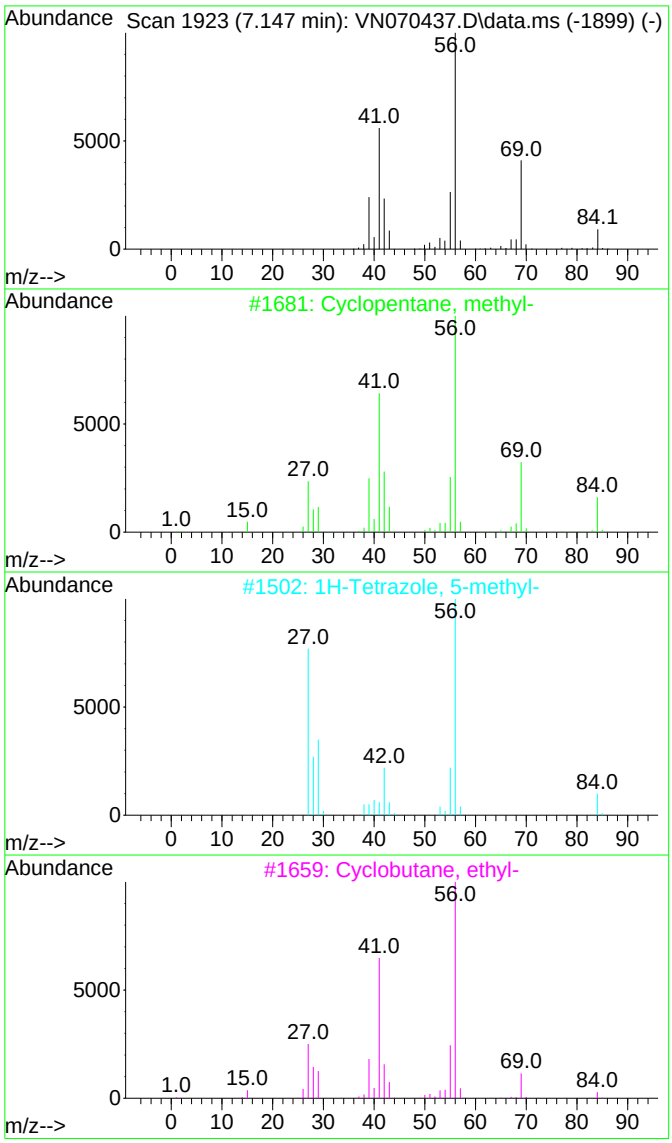
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

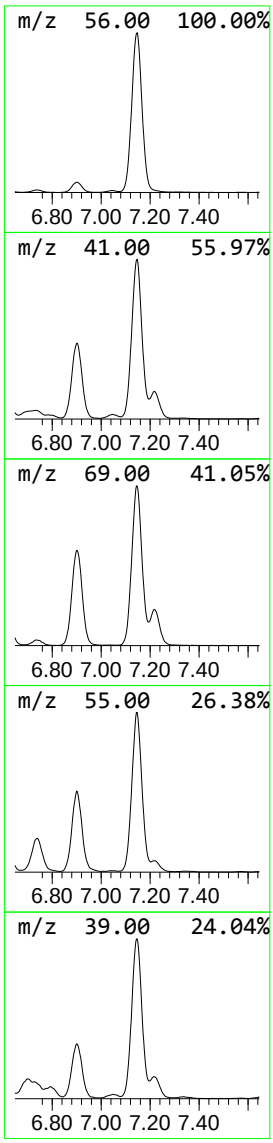
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Cyclopentane, methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.147	81.05 ug/l	11094900	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64
4	Cyclobutane, butyl-	112	C8H16	013152-44-8	64
5	Cyclohexane	84	C6H12	000110-82-7	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

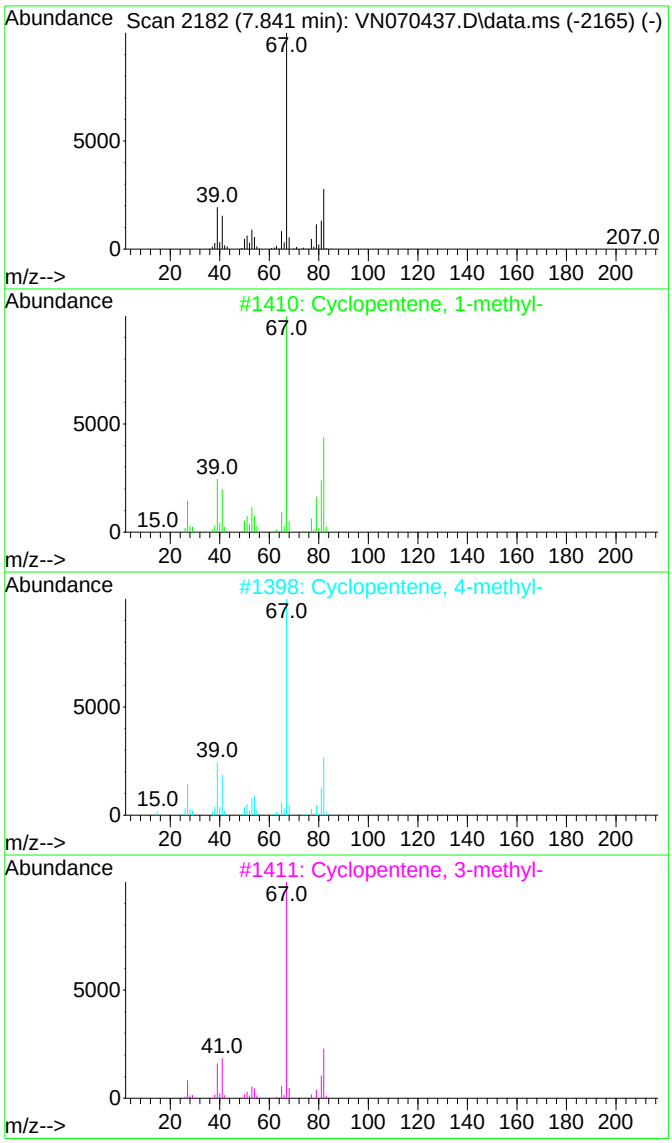
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

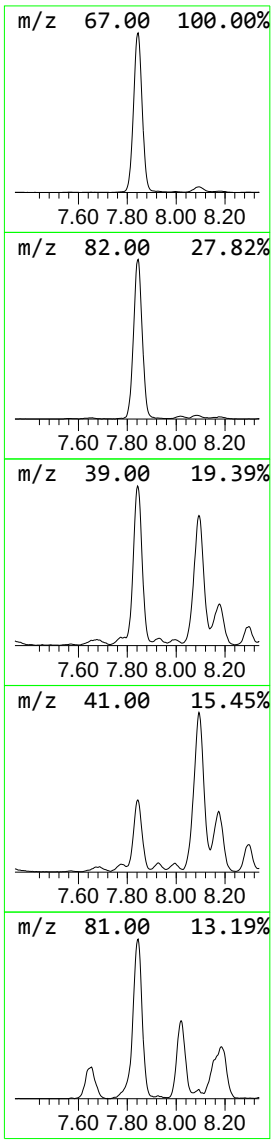
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Cyclopentene, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.841	25.11 ug/l	3437950	Pentafluorobenzene	8.086

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1-methyl-	82	C6H10	000693-89-0	90
2	Cyclopentene, 4-methyl-	82	C6H10	001759-81-5	90
3	Cyclopentene, 3-methyl-	82	C6H10	001120-62-3	86
4	C2H5CH=CHCH=CH2	82	C6H10	000592-48-3	83
5	1,4-Hexadiene, (Z)-	82	C6H10	007318-67-4	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

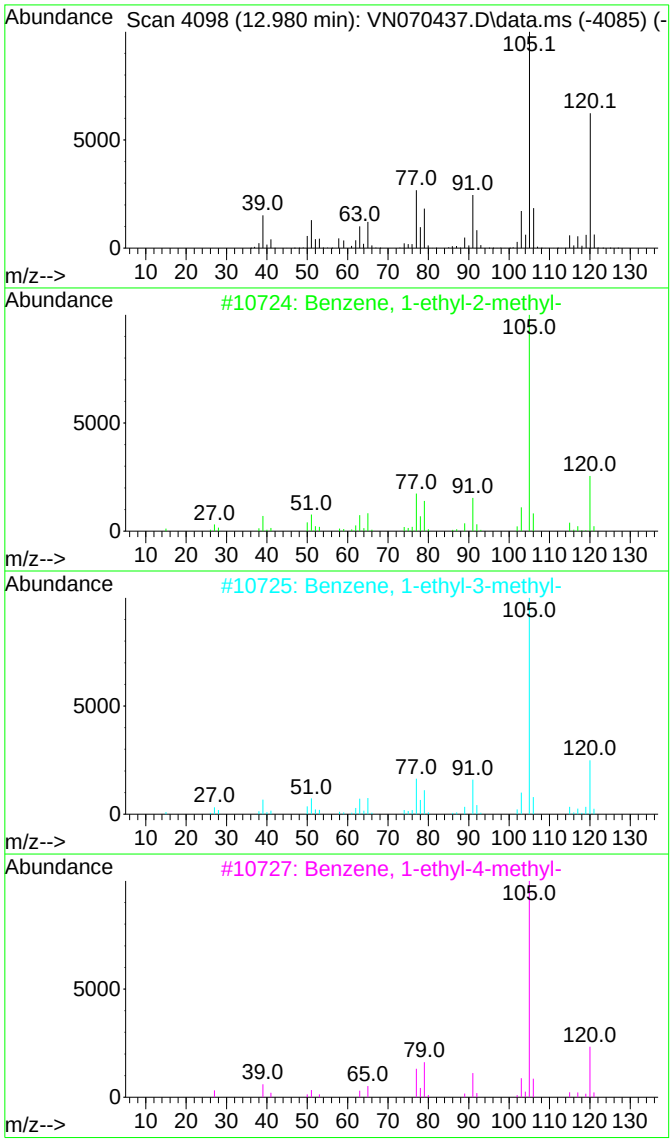
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

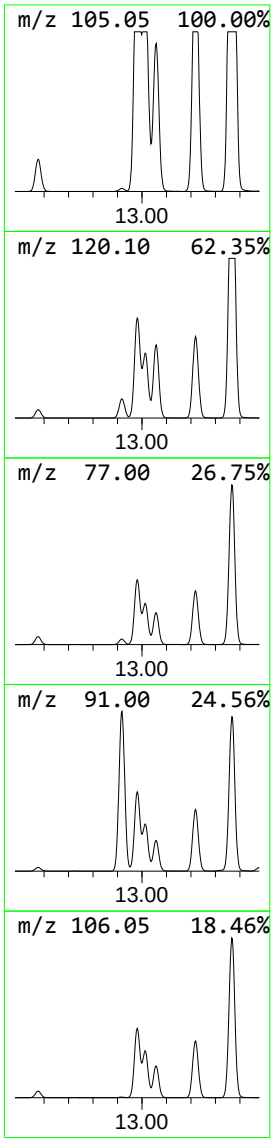
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.980	44.46 ug/l	70217700	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	87
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	87
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	81
4	1,3-Cyclopentadiene, 5-(1-methyl...	120	C9H12	003141-02-4	74
5	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

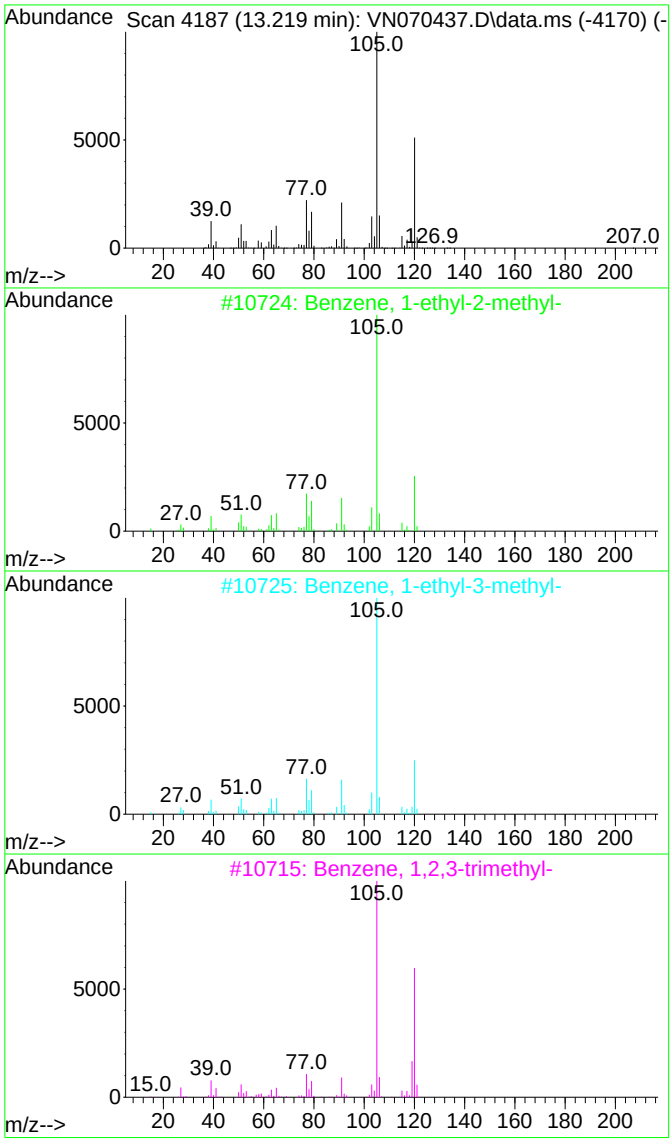
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

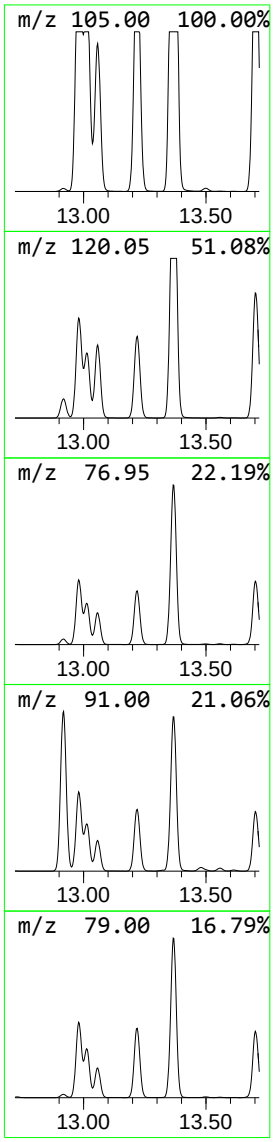
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.219	37.73 ug/l	59591900	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
2	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	90
4	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	87
5	Mesitylene	120	C9H12	000108-67-8	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

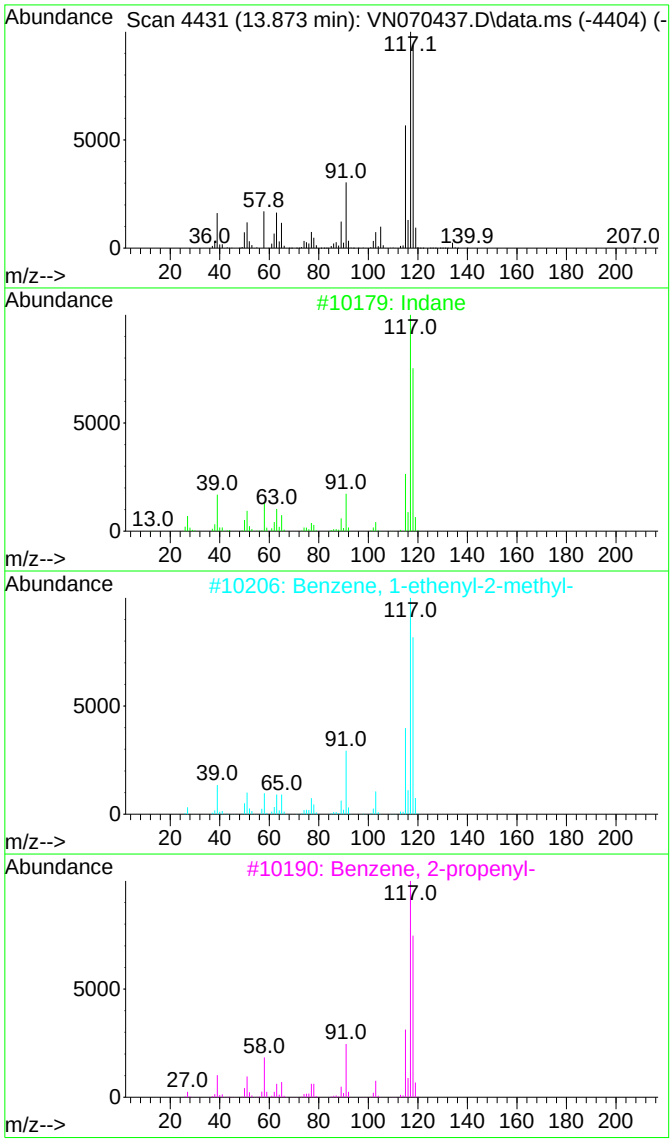
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

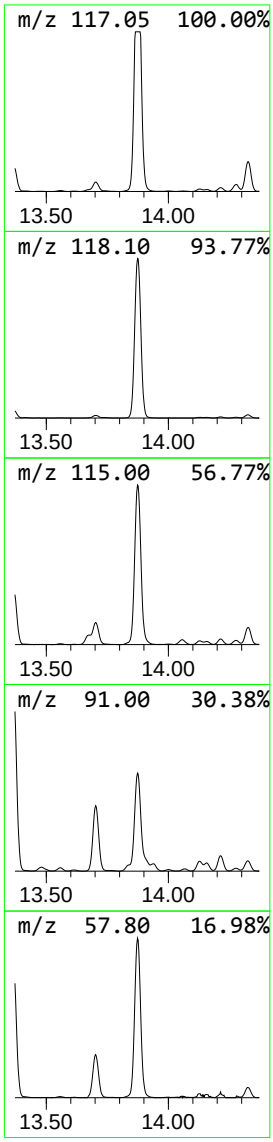
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.873	58.94 ug/l	93096400	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	89
3	Benzene, 2-propenyl-	118	C9H10	000300-57-2	76
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	76
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

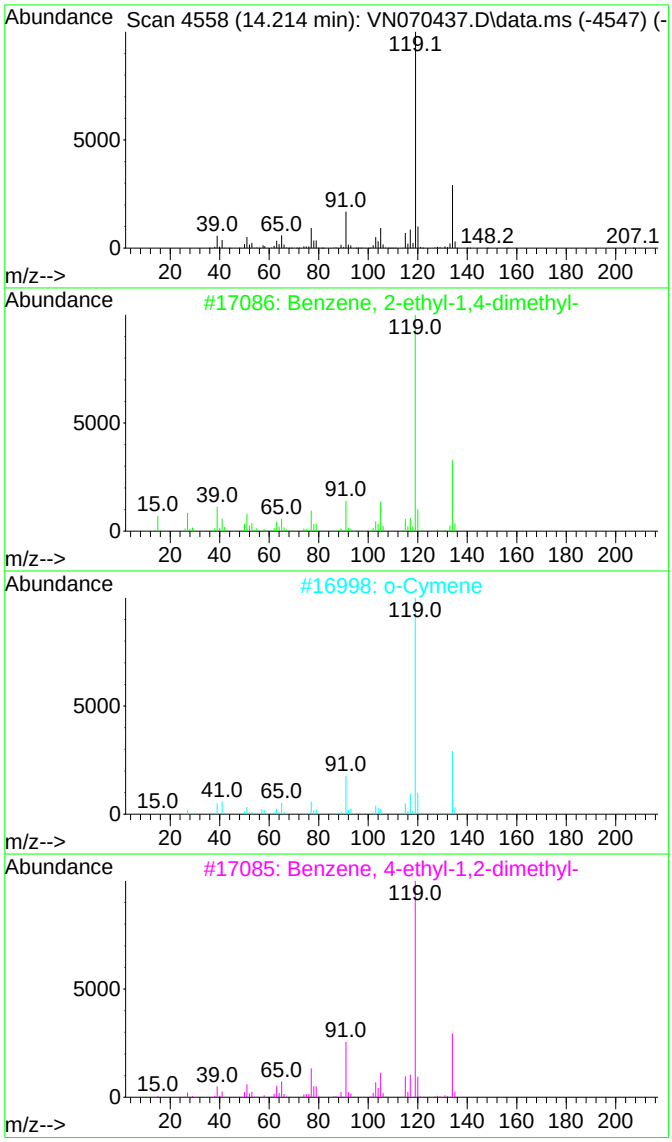
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

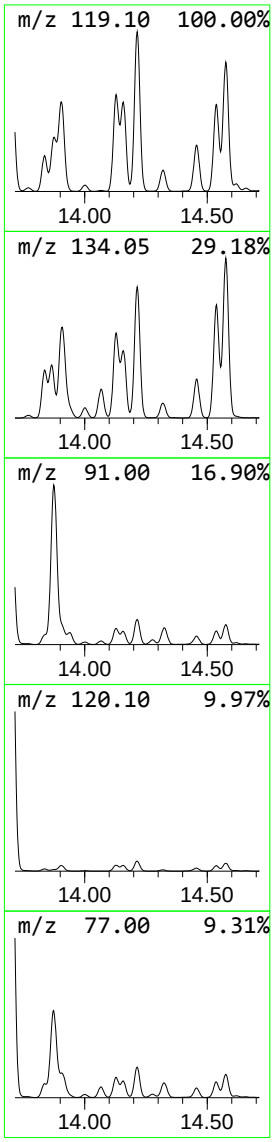
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.214	6.18 ug/l	9753700	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

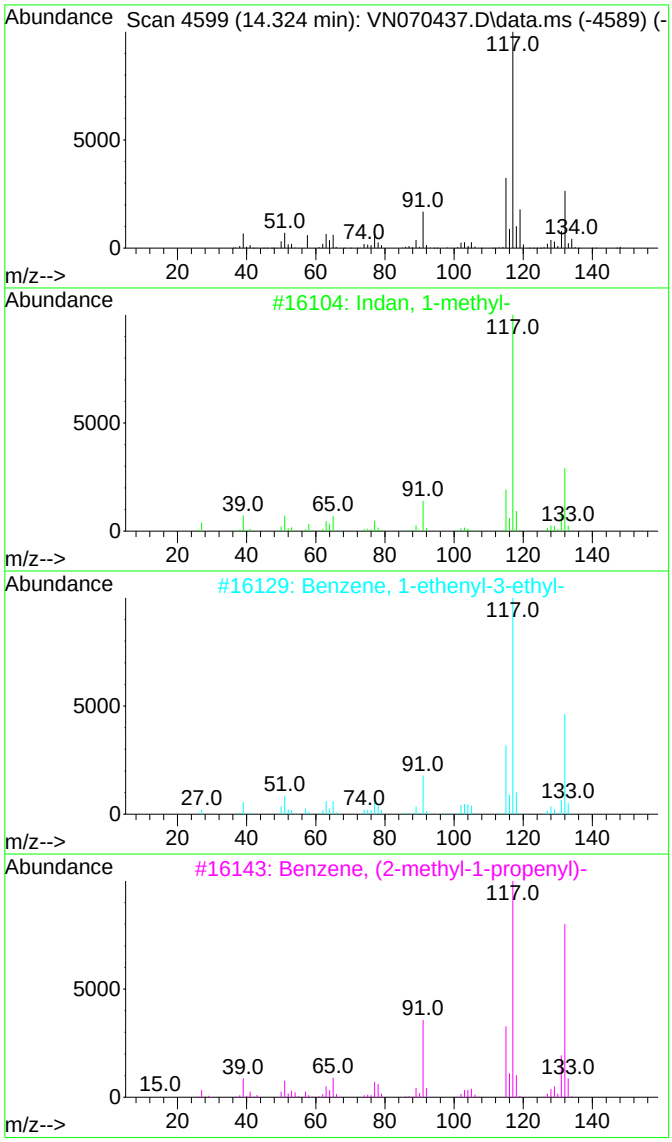
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

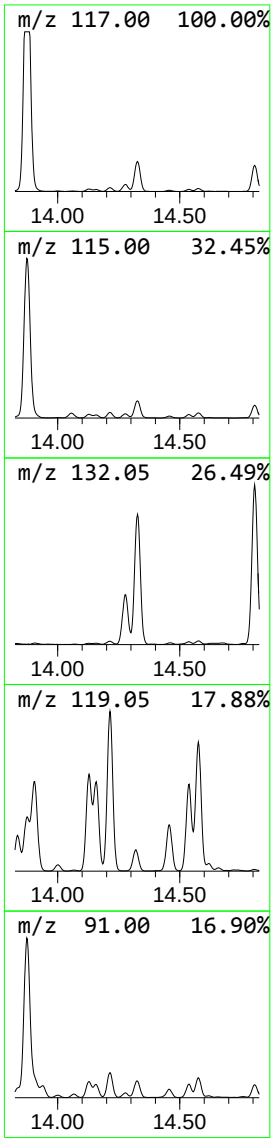
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Indan, 1-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.324	5.52 ug/l	8716060	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	93
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	Benzene, (2-methyl-1-propenyl)-	132	C10H12	000768-49-0	81
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	81
5	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

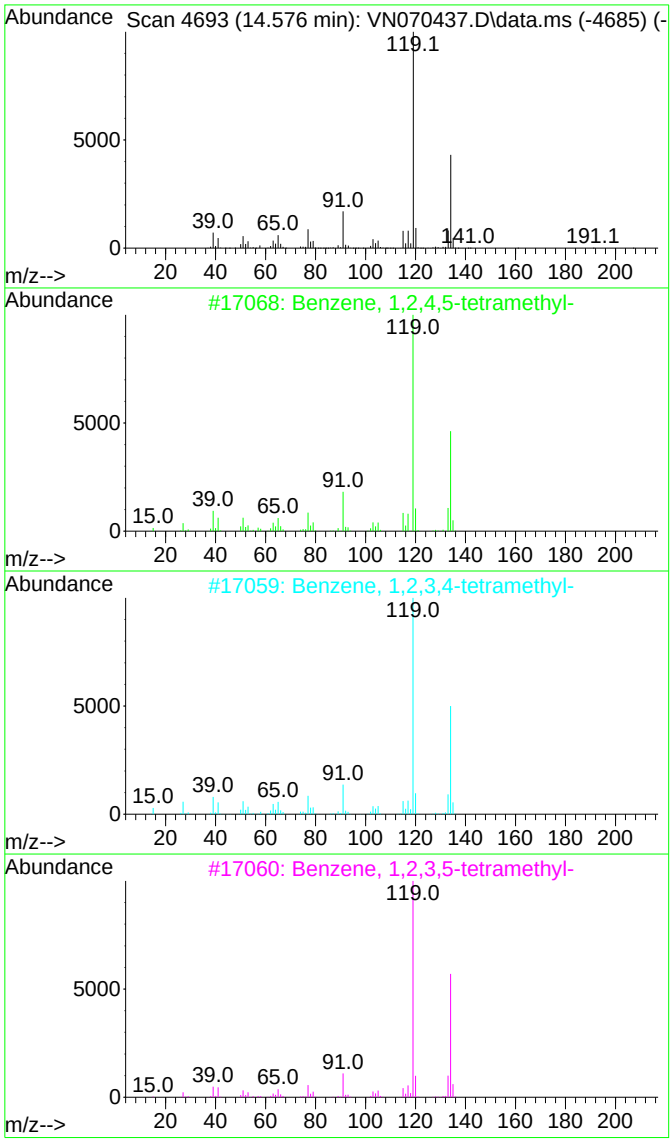
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

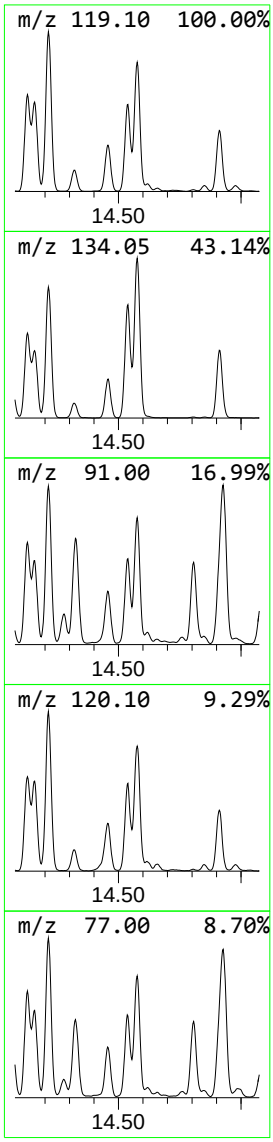
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.576	5.07 ug/l	8004050	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
3	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

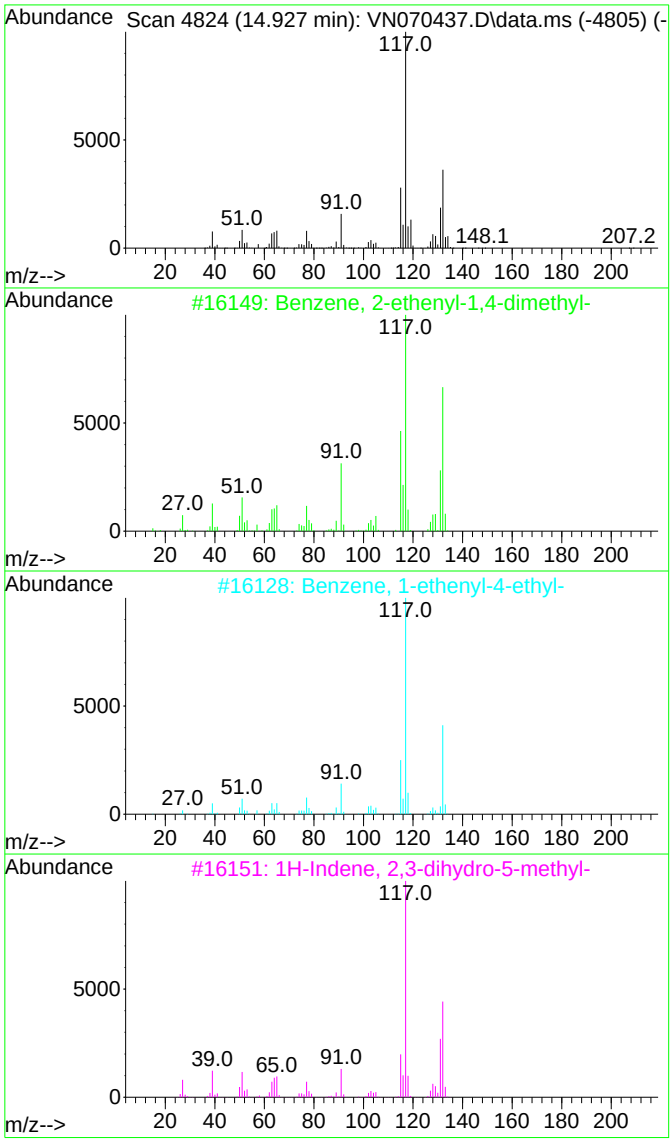
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

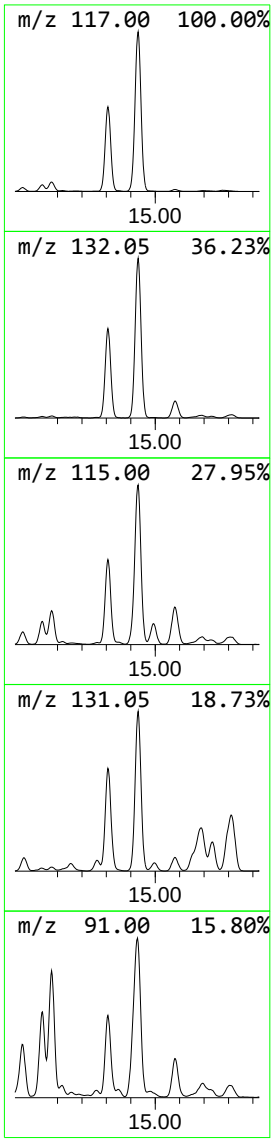
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.927	11.59 ug/l	18311200	1,4-Dichlorobenzene-d4	13.672

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	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	94
4	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	93
5	1-Phenyl-1-butene	132	C10H12	000824-90-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

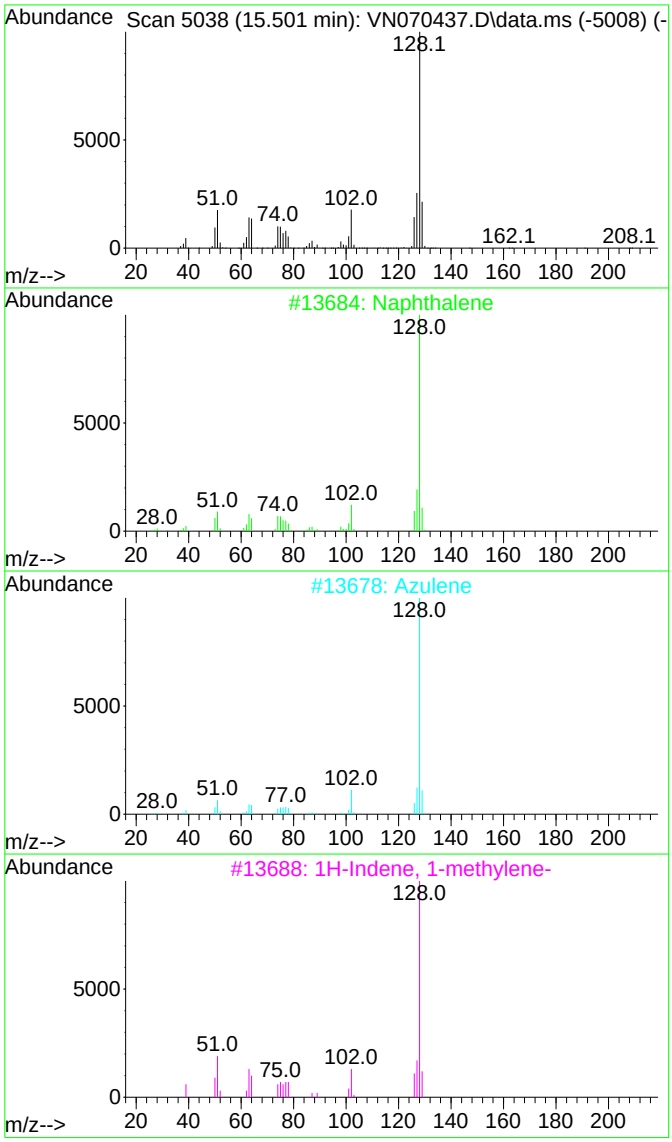
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

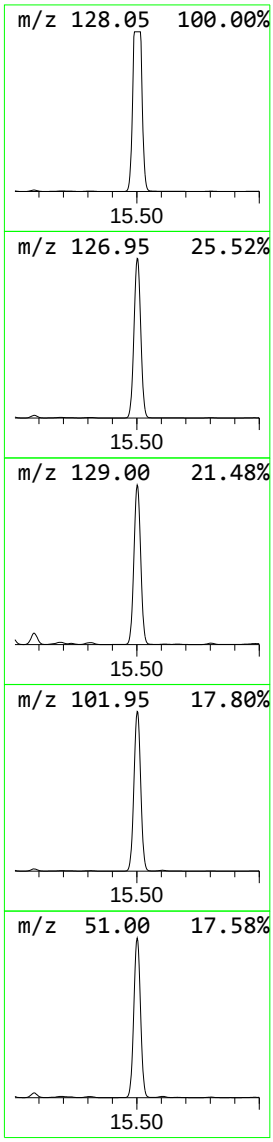
****

Peak Number 14 Azulene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.501	40.55 ug/l	64048900	1,4-Dichlorobenzene-d4	13.672

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene	128	C10H8	000091-20-3	87
2	Azulene	128	C10H8	000275-51-4	81
3	1H-Indene, 1-methylene-	128	C10H8	002471-84-3	81
4	[4.2.2]Propella-2,4,7,9-tetraene	128	C10H8	088090-34-0	62
5	1,4-Benzenedicarbonitrile	128	C8H4N2	000623-26-7	47





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
Data File : VN070437.D
Acq On : 3 Jan 2022 15:10
Operator : JC/MD
Sample : M5251-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 12 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1R

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
Quant Title : SW846 8260

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Sulfur dioxide	2.292	71.5	ug/l	9790720	1	8.086	6844430	50.0
Butane, 2-methyl-	3.140	83.7	ug/l	11463000	1	8.086	6844430	50.0
Pentane	3.513	48.3	ug/l	6613300	1	8.086	6844430	50.0
2-Pentene, 3-me...	6.900	29.0	ug/l	3973020	1	8.086	6844430	50.0
Cyclopentane, m...	7.147	81.0	ug/l	11094900	1	8.086	6844430	50.0
Cyclopentene, 1...	7.841	25.1	ug/l	3437950	1	8.086	6844430	50.0
Benzene, 1-ethy...	12.980	44.5	ug/l	70217700	4	13.672	78975100	50.0
Benzene, 1-ethy...	13.219	37.7	ug/l	59591900	4	13.672	78975100	50.0
Indane	13.873	58.9	ug/l	93096400	4	13.672	78975100	50.0
Benzene, 2-ethy...	14.214	6.2	ug/l	9753700	4	13.672	78975100	50.0
Indan, 1-methyl-	14.324	5.5	ug/l	8716060	4	13.672	78975100	50.0
Benzene, 1,2,4,...	14.576	5.1	ug/l	8004050	4	13.672	78975100	50.0
Benzene, 2-ethe...	14.927	11.6	ug/l	18311200	4	13.672	78975100	50.0
Azulene	15.501	40.5	ug/l	64048900	4	13.672	78975100	50.0