LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

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\SFAMVTR062322WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV026554.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.262	53	69	79	rBV	17049	57283	1.25%	0.185%
2	1.320	81	87	104	rVB	113620	131360	2.87%	0.425%
3	1.449	112	127	152	rBV2	73468	305652	6.69%	0.989%
4	1.581	162	168	181	rVB	100857	132056	2.89%	0.427%
5	2.121	327	336	355	rVB	198332	297845	6.52%	0.964%
6	3.777	830	851	874	rBV2	140348	356208	7.79%	1.153%
7	3.905	879	891	925	rBV2	68640	222299	4.86%	0.719%
8	4.362	1017	1033	1064	rBV2	126242	337571	7.38%	1.093%
9	4.690	1117	1135	1150	rBV2	210308	521452	11.41%	1.688%
10	4.777	1150	1162	1186	rVB4	38781	103113	2.26%	0.334%
11	4.957	1191	1218	1233	rBV5	39729	131440	2.88%	0.425%
12	5.056	1234	1249	1276	rVV2	293153	739506	16.18%	2.393%
13	5.185	1276	1289	1301	rVV4	81823	197830	4.33%	0.640%
14	5.259	1301	1312	1323	rVV4	79654	186742	4.08%	0.604%
15	5.333	1323	1335	1357	rVB3	101003	235689	5.16%	0.763%
16	5.625	1414	1426	1457	rBV	286219	585509	12.81%	1.895%
17	6.076	1551	1566	1573	rBV	165383	334010	7.31%	1.081%
18	6.133	1575	1584	1605	rVB3	379110	843381	18.45%	2.729%
19	6.368	1645	1657	1673	rVB2	41181	79260	1.73%	0.257%
20	6.992	1839	1851	1876	rBV	73904	146317	3.20%	0.474%
21	7.268	1924	1937	1943	rBV3	34942	66547	1.46%	0.215%
22	7.320	1943	1953	1970	rVB	362463	640336	14.01%	2.072%
23	7.629	2031	2049	2053	rBV	37012	60205	1.32%	0.195%
24	7.657	2054	2058	2073	rVB3	46423	73567	1.61%	0.238%
25	8.092	2182	2193	2224	rBV	248049	467170	10.22%	1.512%
26	8.291	2246	2255	2265	rVB	37453	60789	1.33%	0.197%
27	8.854	2419	2430	2447	rBV	393849	641207	14.03%	2.075%
28	9.014	2469	2480	2493	rBV	142984	228256	4.99%	0.739%
29	9.146	2513	2521	2540	rVB	54318	90611	1.98%	0.293%
30	9.931	2751	2765	2791	rBV	1711654	2554688	55.88%	8.268%
31	10.217	2841	2854	2875	rBV	109581	171852	3.76%	0.556%
32	10.352	2882	2896	2923	rBV	3085919	4571515	100.00%	14.795%
33	10.915	3061	3071	3088	rVB	298010	445959	9.76%	1.443%
34	11.056	3107	3115	3119	rBV	37353	52763	1.15%	0.171%
35	11.088	3119	3125	3141	rVB	61647	97528	2.13%	0.316%
36	11.249	3162	3175	3191	rBV	417706	639939	14.00%	2.071%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

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\SFAMVTR062322WMA.M
Title : TRACE VOA SFAM1.0

37	11.336	3191	3202	3217	rVB	171384	262755	5.75%	0.850%
38	11.525	3249	3261	3280	rBV2	1200781	2140146	46.81%	6.926%
39	11.625	3280	3292	3318	rVV4	461685	892445	19.52%	2.888%
40	11.744	3318	3329	3342	rVV2	75600	116974	2.56%	0.379%
41	11.818	3342	3352	3364	rVB	32834	50999	1.12%	0.165%
42	11.911	3370	3381	3386	rBV	59568	92040	2.01%	0.298%
43	12.014	3399	3413	3434	rBV	1511642	2293108	50.16%	7.421%
44	12.114	3434	3444	3466	rVV3	247136	596186	13.04%	1.929%
45	12.313	3492	3506	3517	rVB3	35853	63056	1.38%	0.204%
46	12.410	3517	3536	3546	rBV	1083666	1542364	33.74%	4.992%
47	12.464	3546	3553	3570	rVV	230704	350223	7.66%	1.133%
48	12.548	3570	3579	3589	rVV	39852	56617	1.24%	0.183%
49	12.722	3622	3633	3649	rVB	457186	686329	15.01%	2.221%
50	12.879	3663	3682	3693	rBV2	1745704	2691054	58.87%	8.709%
51	12.998	3711	3719	3726	rBV	52414	76548	1.67%	0.248%
52	13.050	3727	3735	3756	rVB6	60162	119037	2.60%	0.385%
53	13.268	3794	3803	3818	rBV2	93304	154477	3.38%	0.500%
54	13.400	3837	3844	3855	rVB	53622	81297	1.78%	0.263%
55	13.500	3862	3875	3887	rBV	988752	1489809	32.59%	4.822%
56	16.483	4791	4803	4821	rBV	207658	335968	7.35%	1.087%

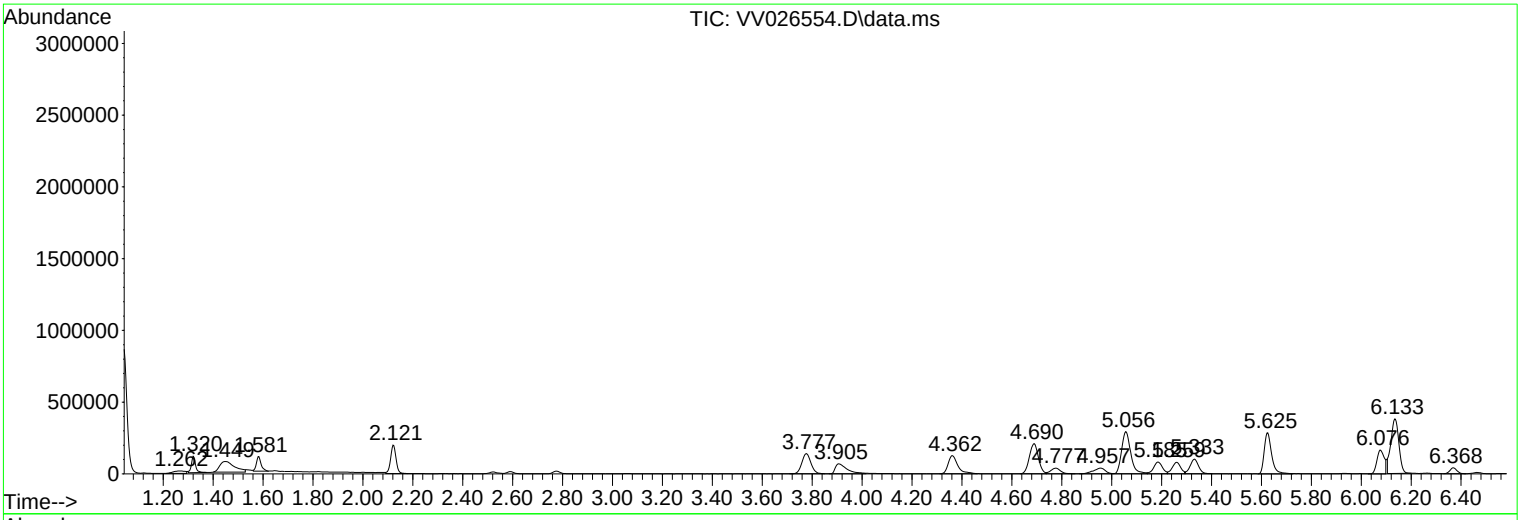
Sum of corrected areas: 30898887

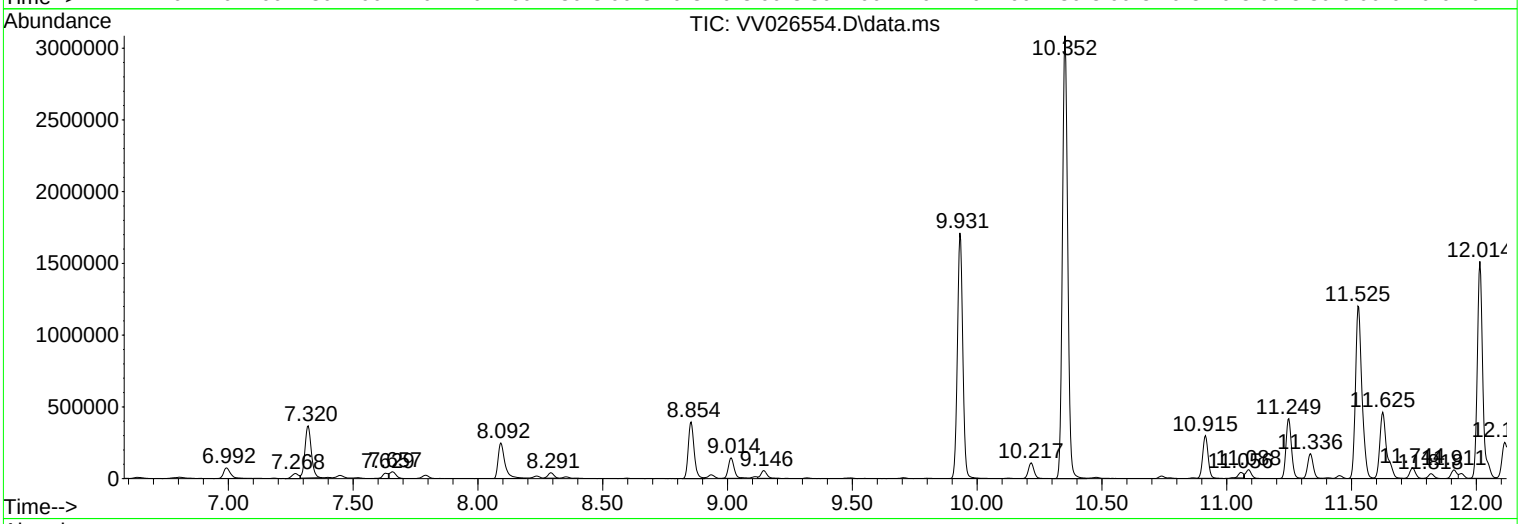
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

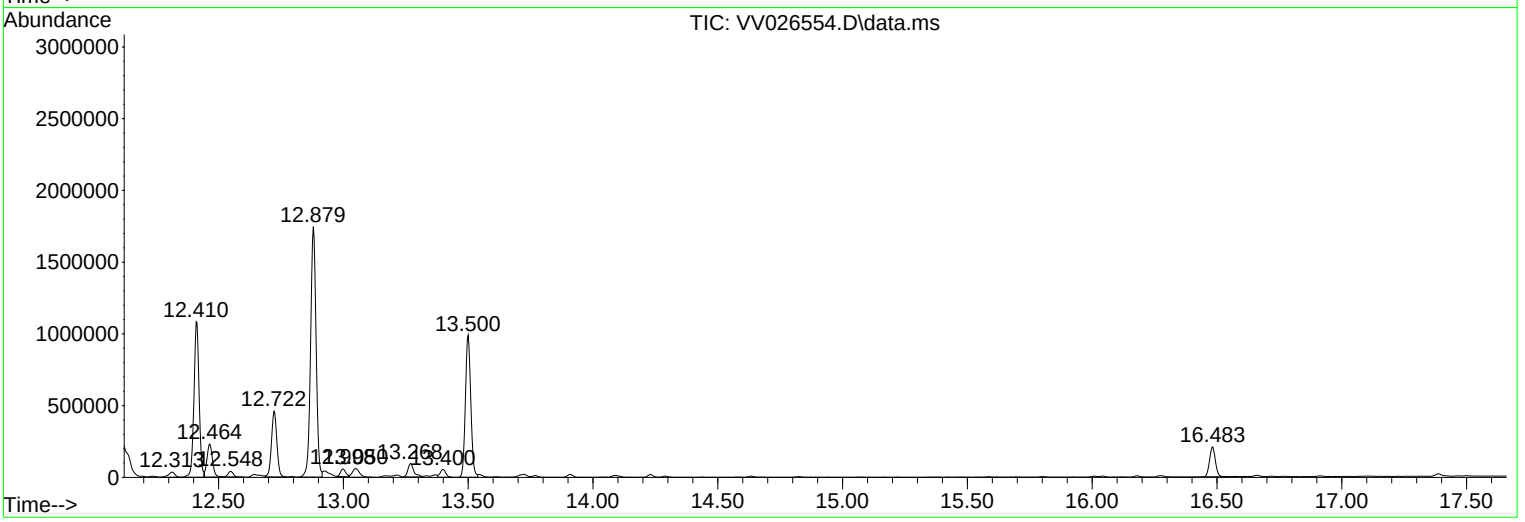
Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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







Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

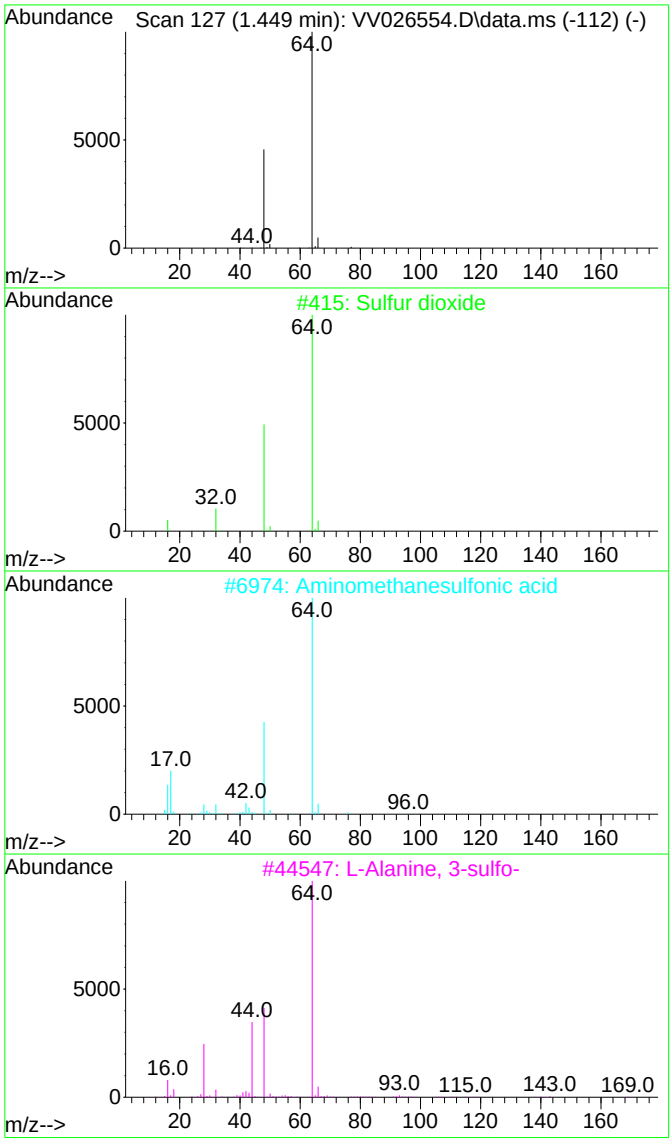
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

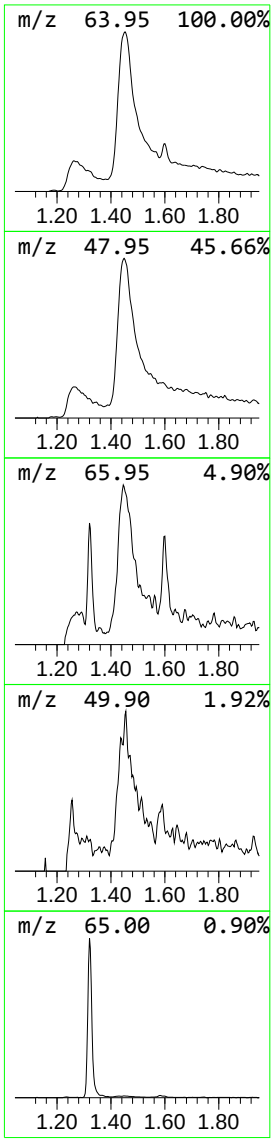
****

Peak Number 1 Sulfur dioxide Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.449	2.61 ug/L	305652	1,4-Difluorobenzene	5.625

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	Sulfur dioxide	64	O2S	007446-09-5	5
5	Ethene, 1,1-difluoro-	64	C2H2F2	000075-38-7	4





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

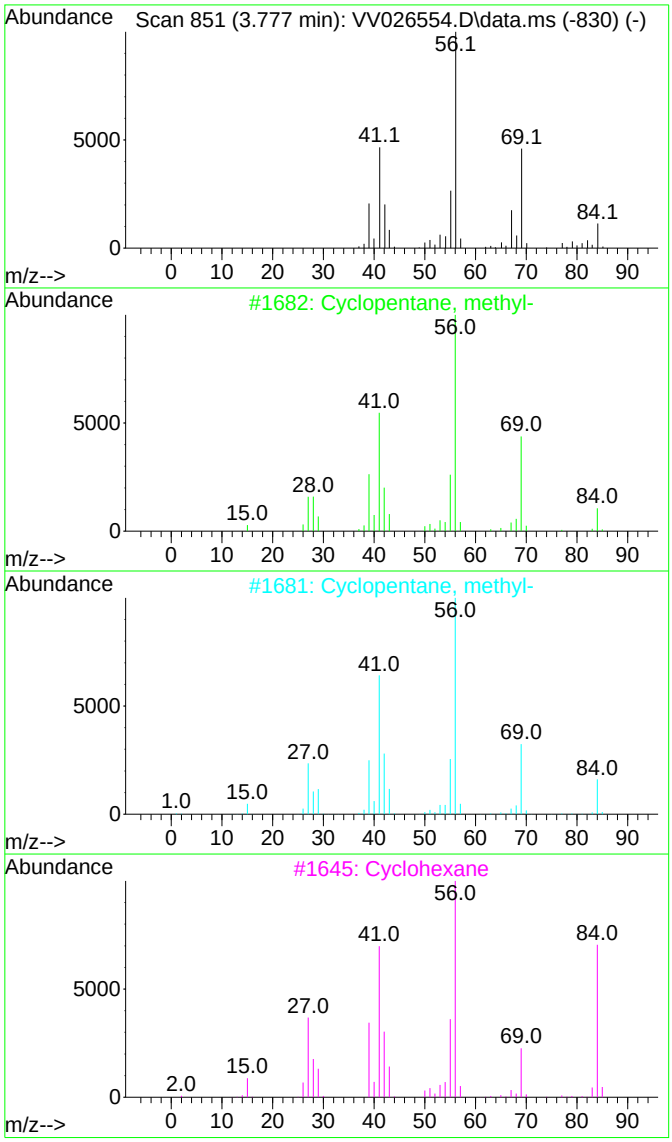
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

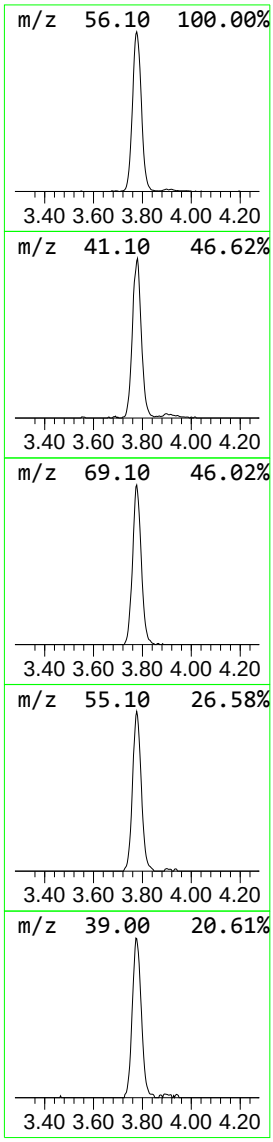
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Cyclic3.777 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.777	3.04 ug/L	356208	1,4-Difluorobenzene	5.625

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	83
3	Cyclohexane	84	C6H12	000110-82-7	74
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	72
5	1-Pentene, 2-methyl-	84	C6H12	000763-29-1	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

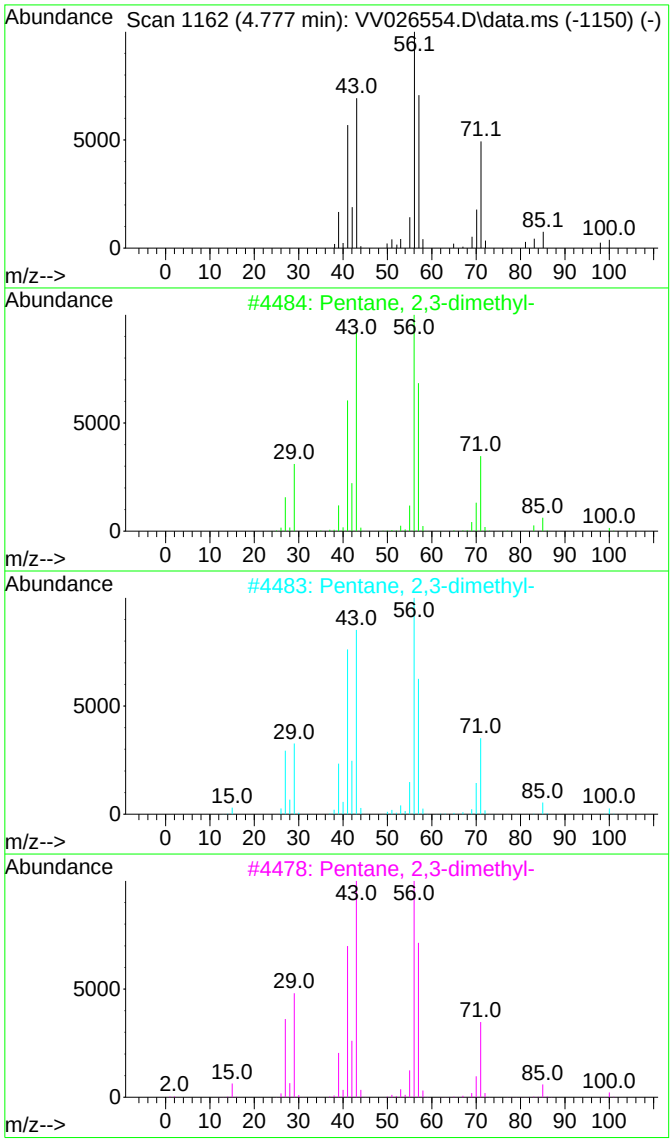
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

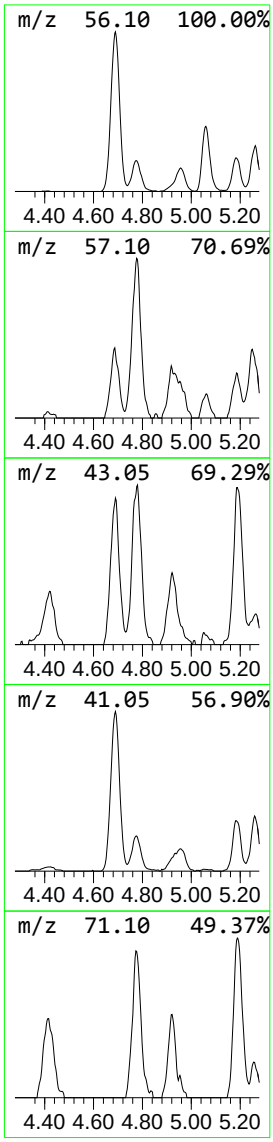
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.777	0.88 ug/L	103113	1,4-Difluorobenzene	5.625

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	91
3	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	90
4	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	87
5	Hexane, 2,2,4-trimethyl-	128	C9H20	016747-26-5	50





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

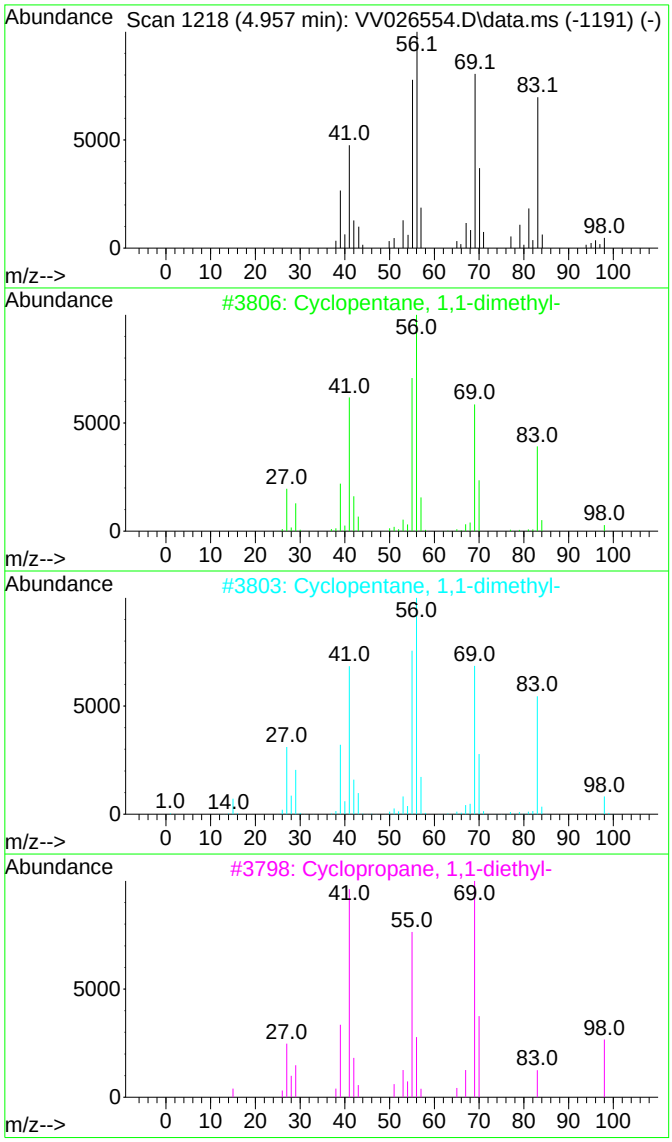
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

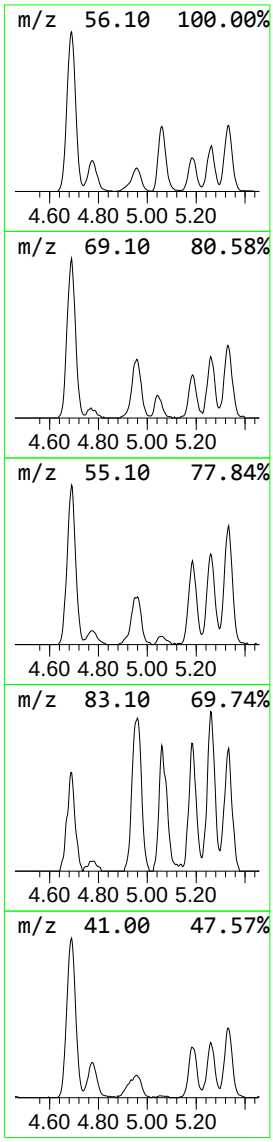
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Cyclic4.957 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.957	1.12 ug/L	131440	1,4-Difluorobenzene	5.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	86
2	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	78
3	Cyclopropane, 1,1-diethyl-	98	C7H14	001003-19-6	55
4	2-Hexene, 3-methyl-, (Z)-	98	C7H14	010574-36-4	38
5	3-Hexene, 2-methyl-, (Z)-	98	C7H14	015840-60-5	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

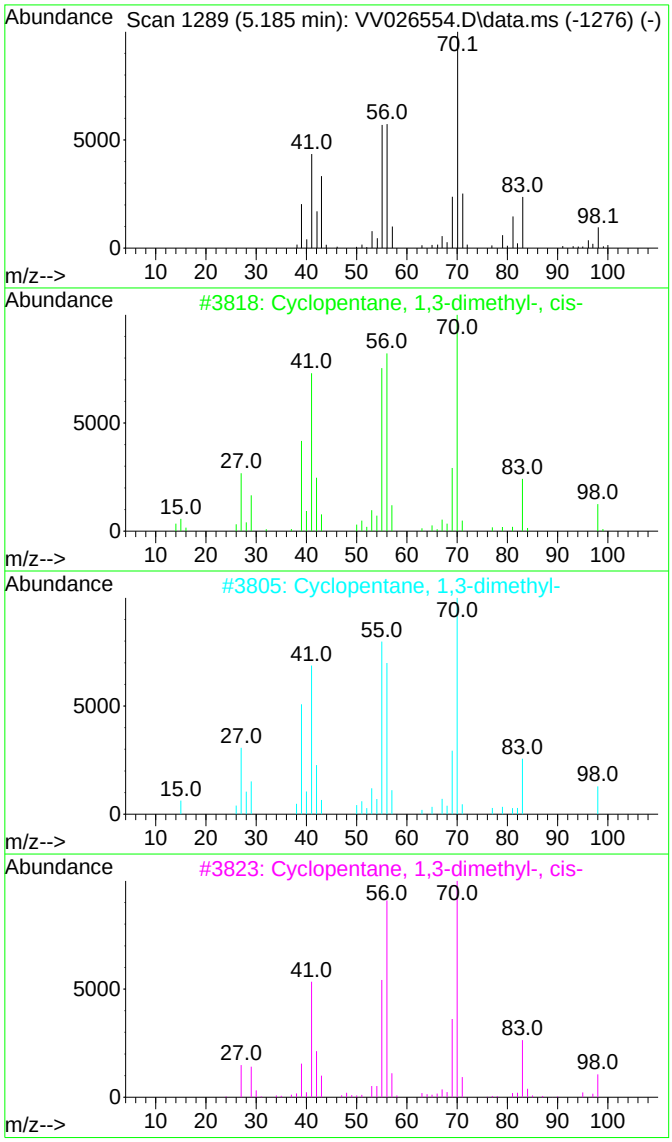
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

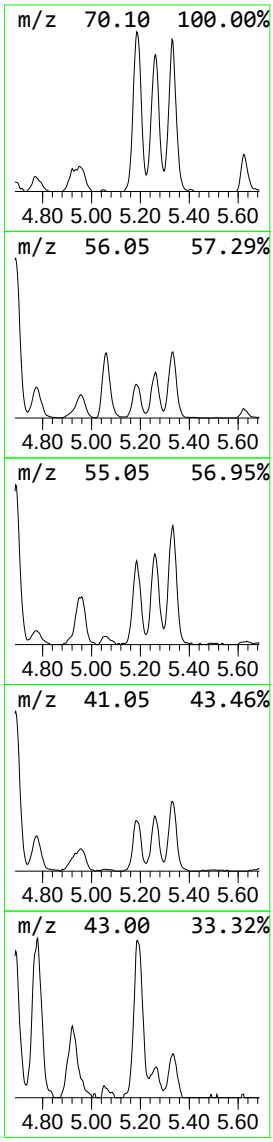
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Cyclic5.185 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.185	1.69 ug/L	197830	1,4-Difluorobenzene	5.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	87
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	64
3	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	64
4	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	60
5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	53





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

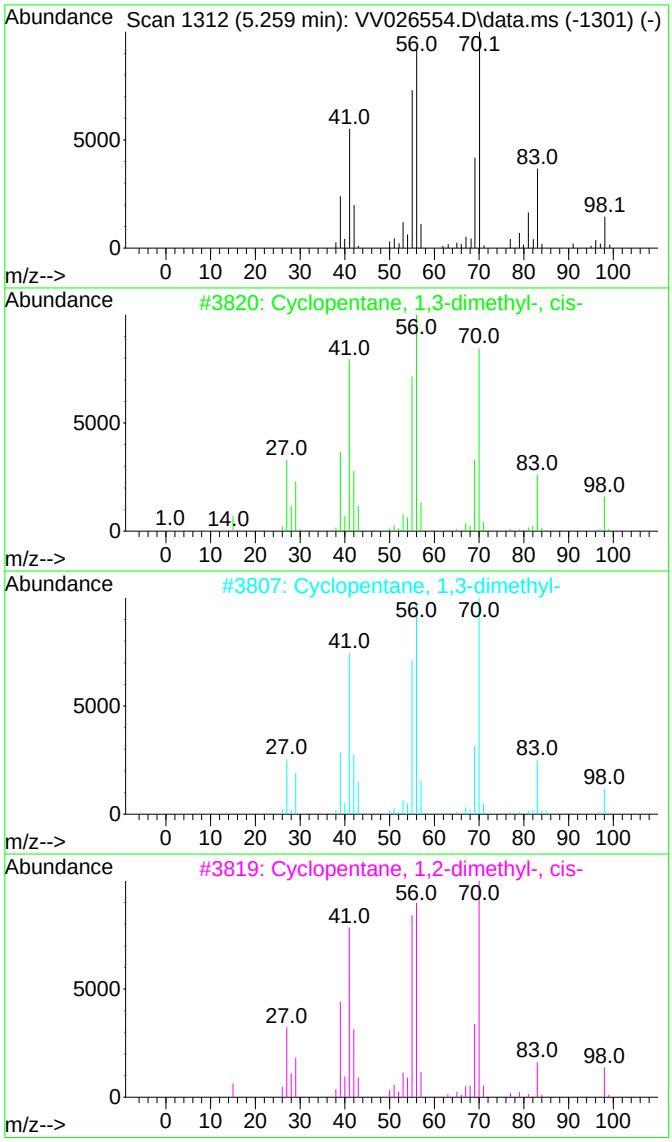
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

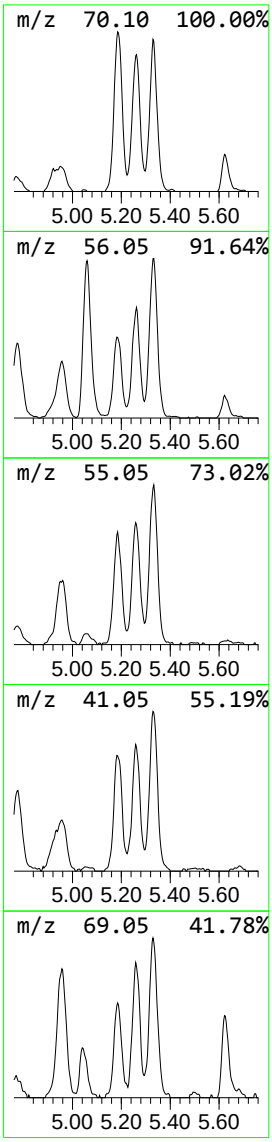
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic5.259 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.259	1.59 ug/L	186742	1,4-Difluorobenzene	5.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	91
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	90
3	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	90
4	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	87
5	Cyclopentane, 1,2-dimethyl-, cis-	98	C7H14	001192-18-3	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

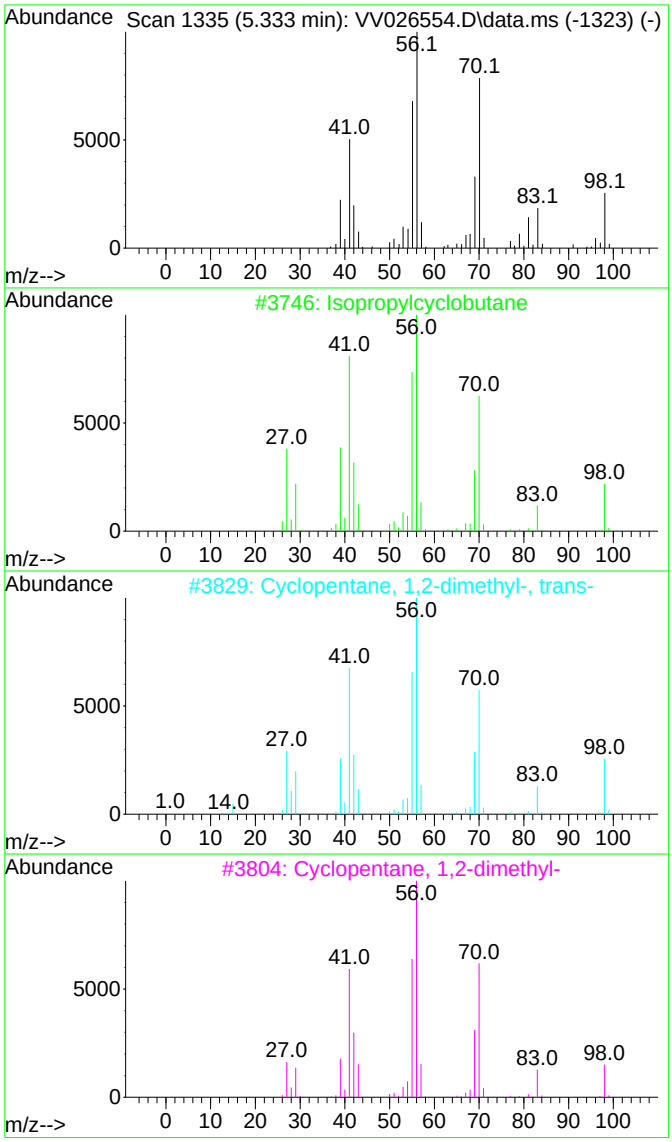
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

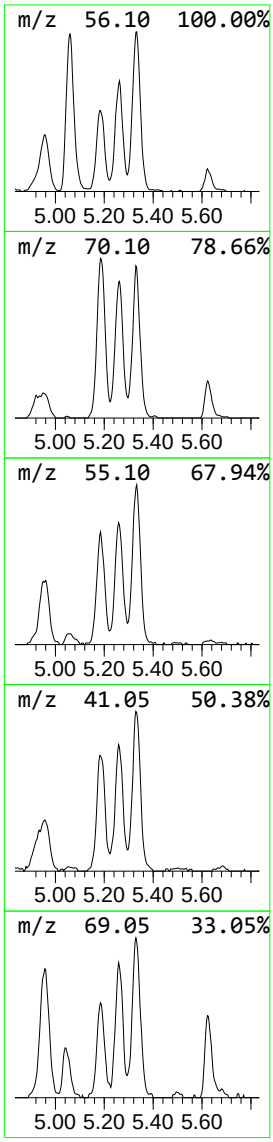
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 (DEL) Alkane: Cyclic5.333 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.333	2.01 ug/L	235689	1,4-Difluorobenzene	5.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Isopropylcyclobutane	98	C7H14	000872-56-0	93
2	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	91
3	Cyclopentane, 1,2-dimethyl-	98	C7H14	002452-99-5	90
4	Cyclopentane, 1,2-dimethyl-, trans-	98	C7H14	000822-50-4	80
5	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

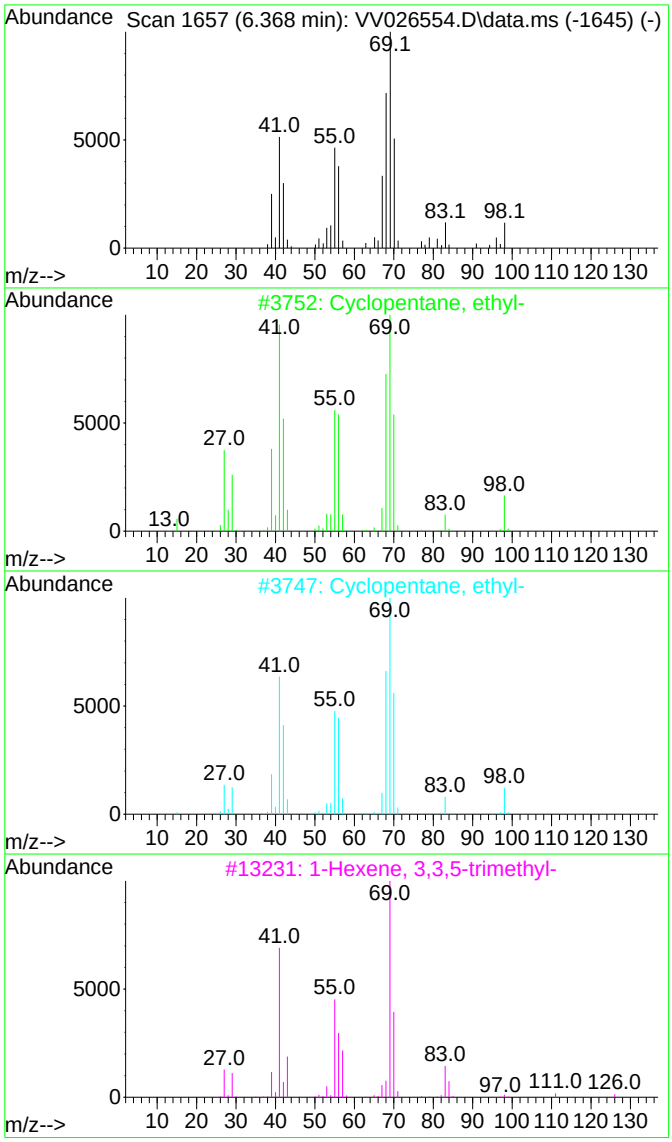
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

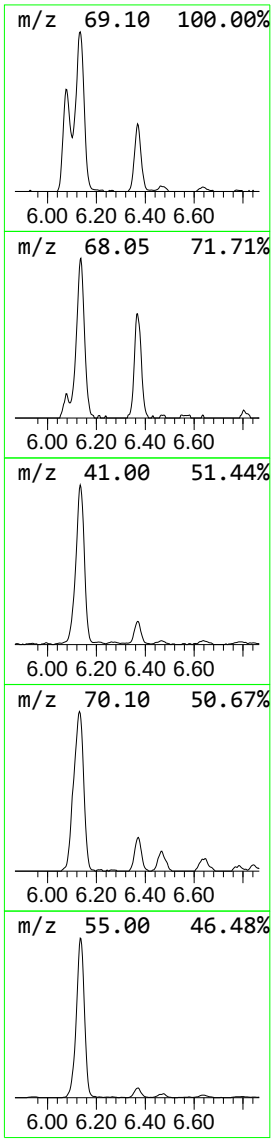
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 (DEL) Alkane: Cyclic6.368 Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.368	0.68 ug/L	79260	1,4-Difluorobenzene	5.625

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, ethyl-	98	C7H14	001640-89-7	90
2	Cyclopentane, ethyl-	98	C7H14	001640-89-7	90
3	1-Hexene, 3,3,5-trimethyl-	126	C9H18	013427-43-5	47
4	1-Pentene, 3-ethyl-	98	C7H14	004038-04-4	47
5	1-Pentene, 2,3-dimethyl-	98	C7H14	003404-72-6	47





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Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

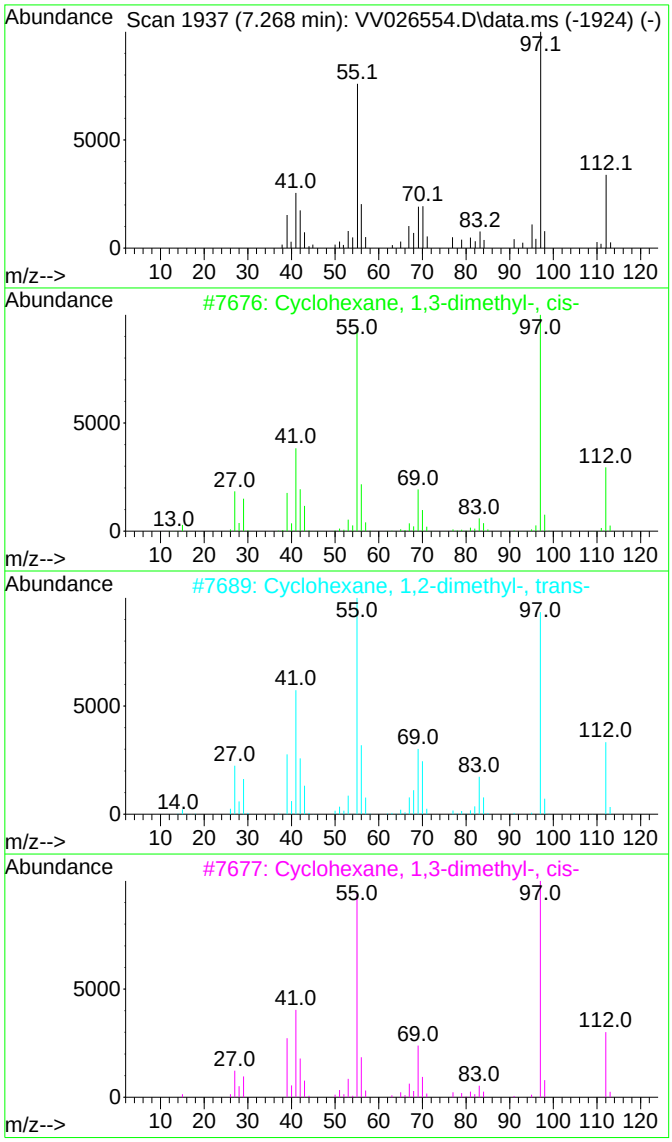
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

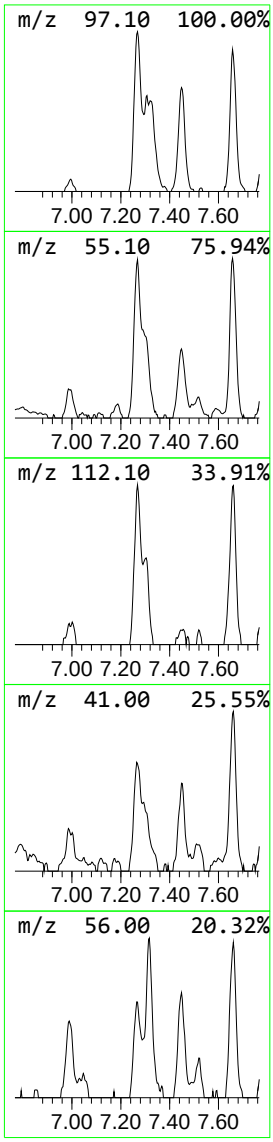
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 (DEL) Alkane: Cyclic7.268 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.268	0.52 ug/L	66547	Chlorobenzene-d5	8.854

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	93
2	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	87
3	Cyclohexane, 1,3-dimethyl-, cis-	112	C8H16	000638-04-0	87
4	Cyclohexane, 1,2-dimethyl-, trans-	112	C8H16	006876-23-9	83
5	Cyclohexane, 1,2-dimethyl- (cis/...	112	C8H16	000583-57-3	83





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Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

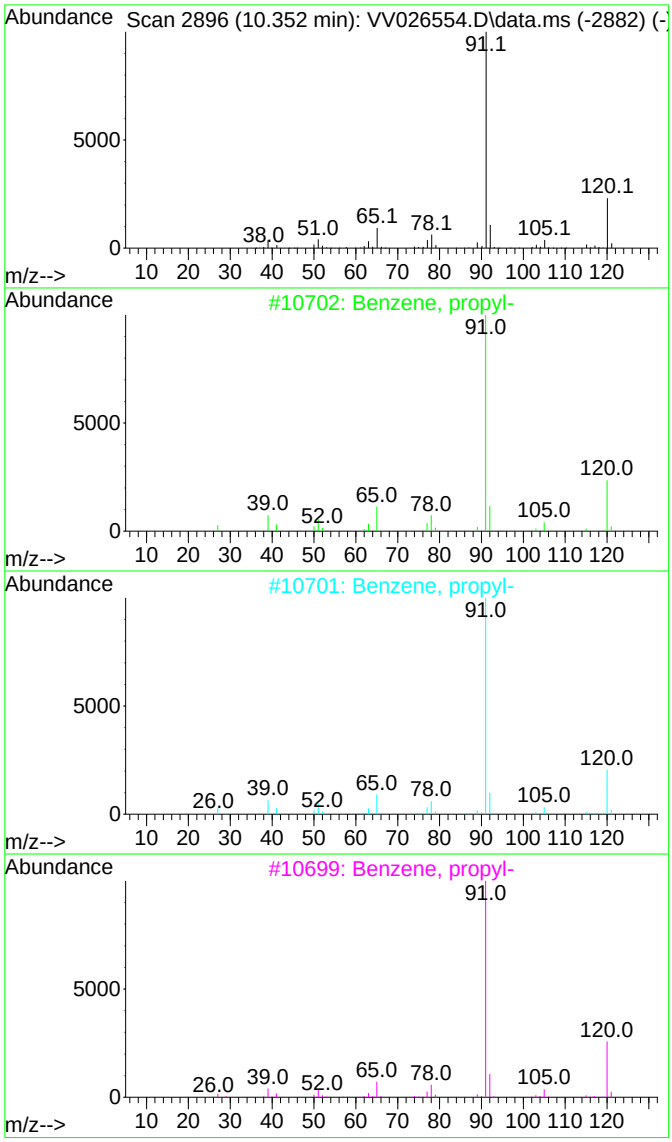
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

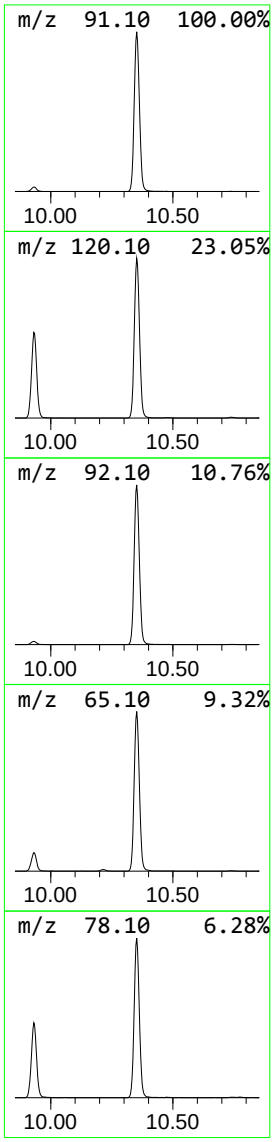
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, propyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.352	35.72 ug/L	4571520	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	93
2	Benzene, propyl-	120	C9H12	000103-65-1	90
3	Benzene, propyl-	120	C9H12	000103-65-1	90
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5	Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

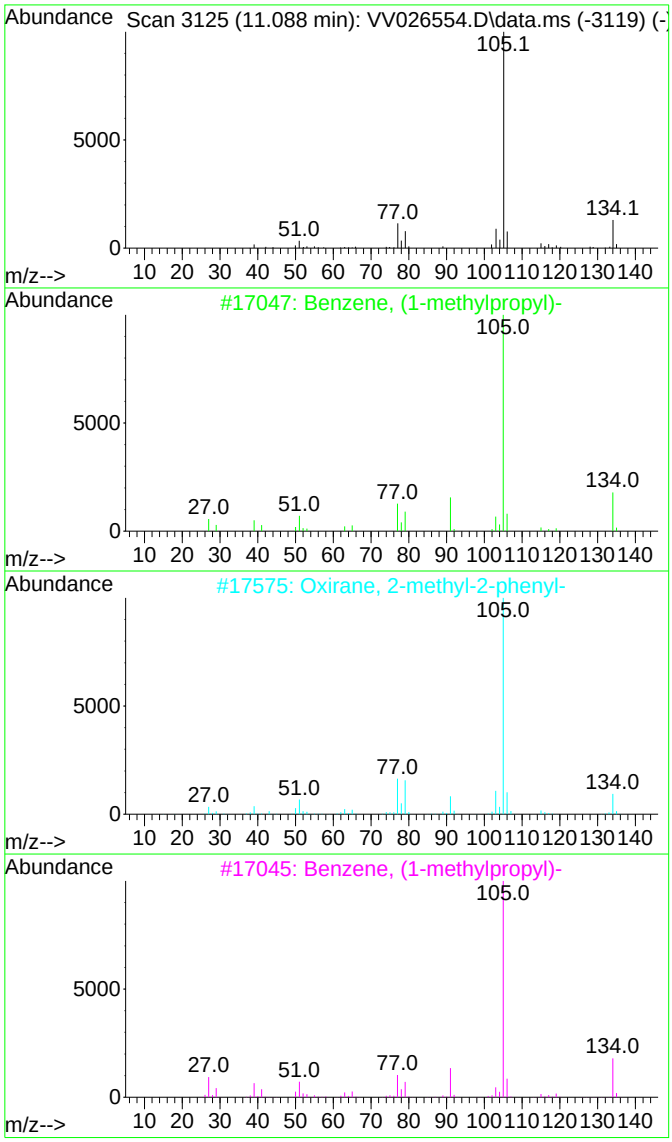
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

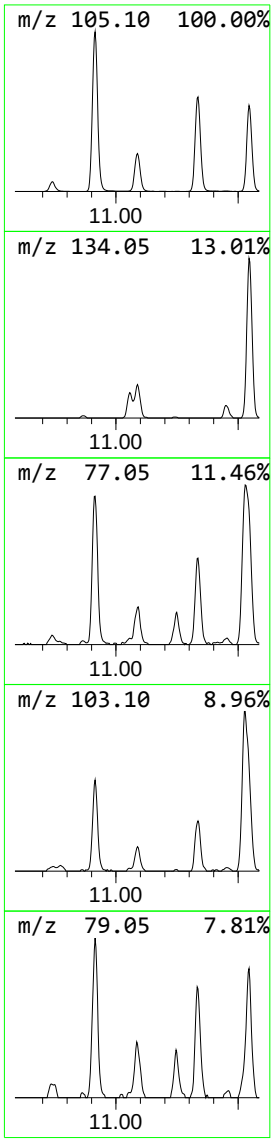
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, (1-methylpropyl)- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.088	0.76 ug/L	97528	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	87
2	Oxirane, 2-methyl-2-phenyl-	134	C9H10O	002085-88-3	87
3	Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	80
4	Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	80
5	2-Tolyloxirane	134	C9H10O	002783-26-8	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

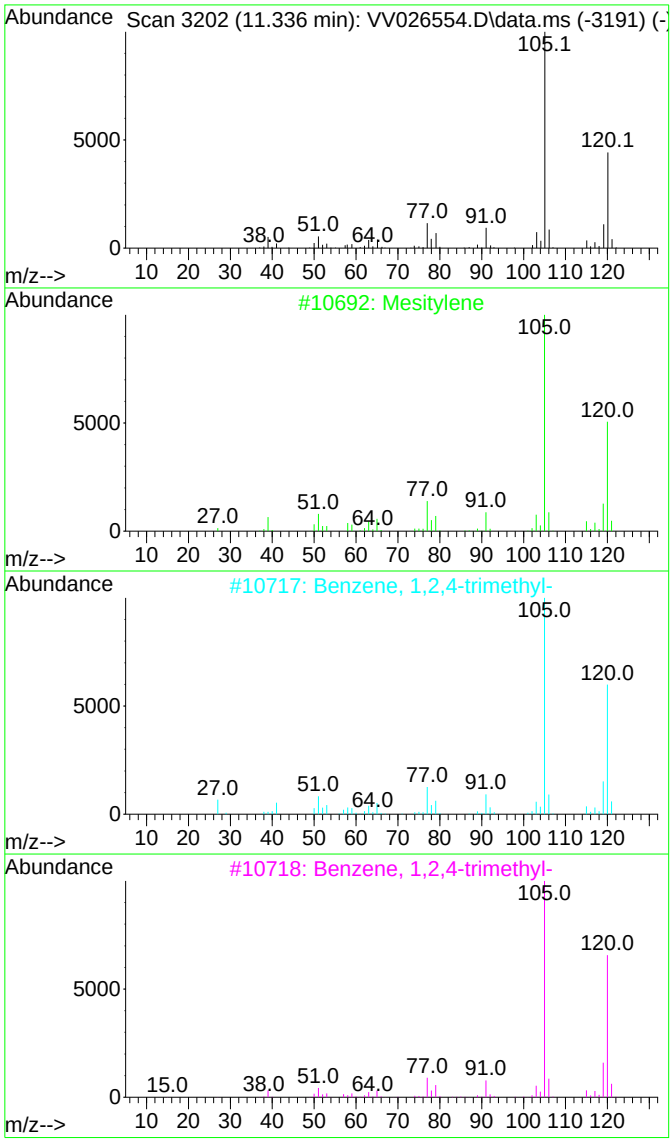
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

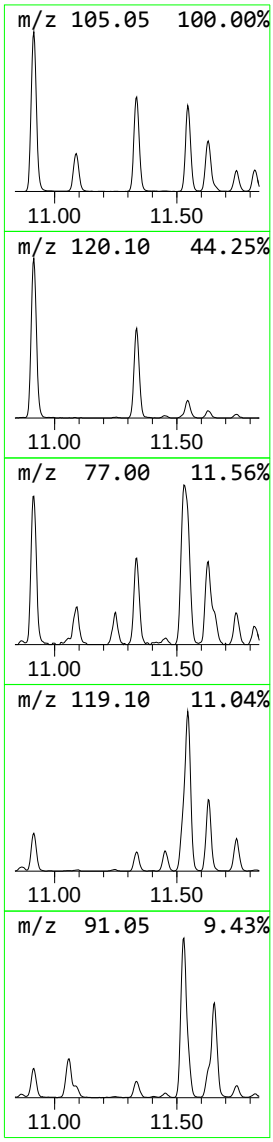
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,2,3-trimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.336	2.05 ug/L	262755	1,4-Dichlorobenzene-d4	11.249

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

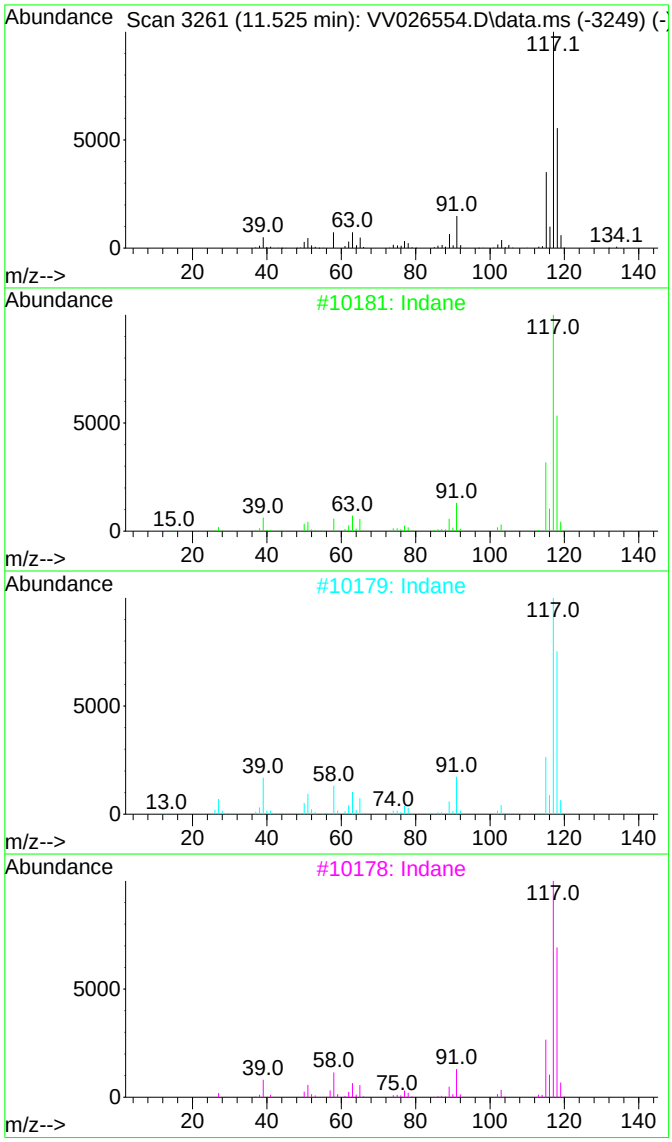
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

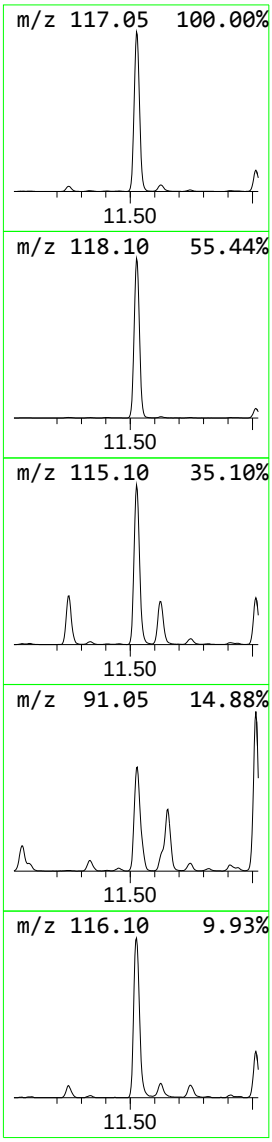
****

Peak Number 14 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.525	16.72 ug/L	2140150	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	95
2	Indane	118	C9H10	000496-11-7	91
3	Indane	118	C9H10	000496-11-7	87
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
5	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

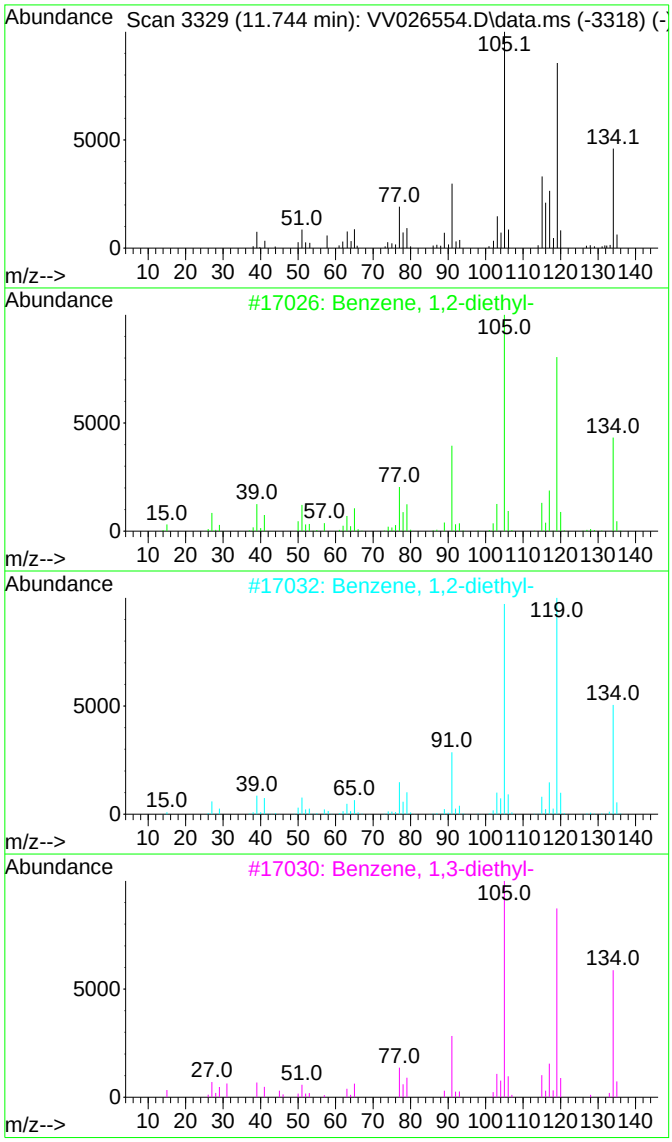
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

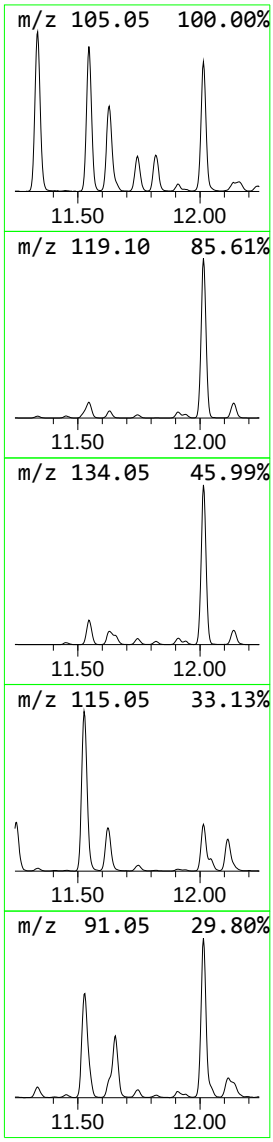
****

Peak Number 15 Benzene, 1,2-diethyl- Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.744	0.91 ug/L	116974	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	87
2	Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	76
3	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76
4	Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	76
5	Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

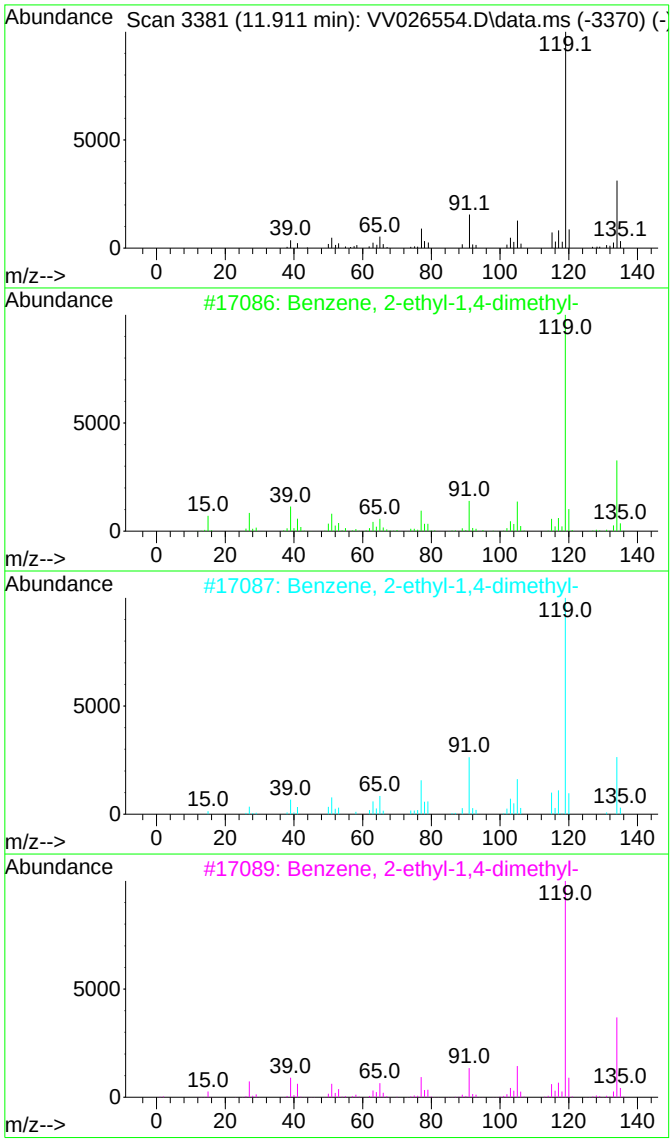
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

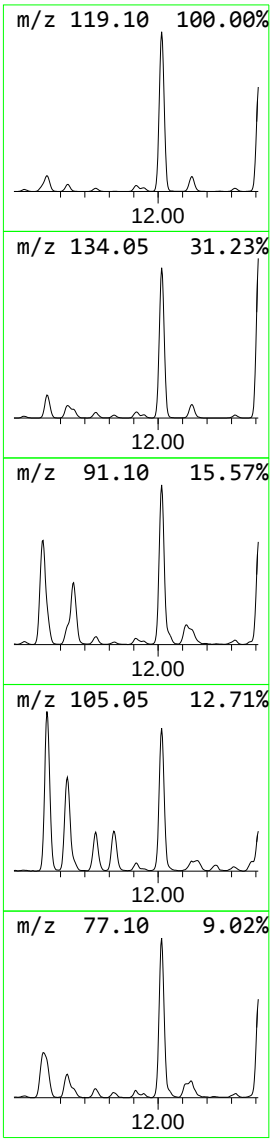
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 16 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.911	0.72 ug/L	92040	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	97
2	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	96
3	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

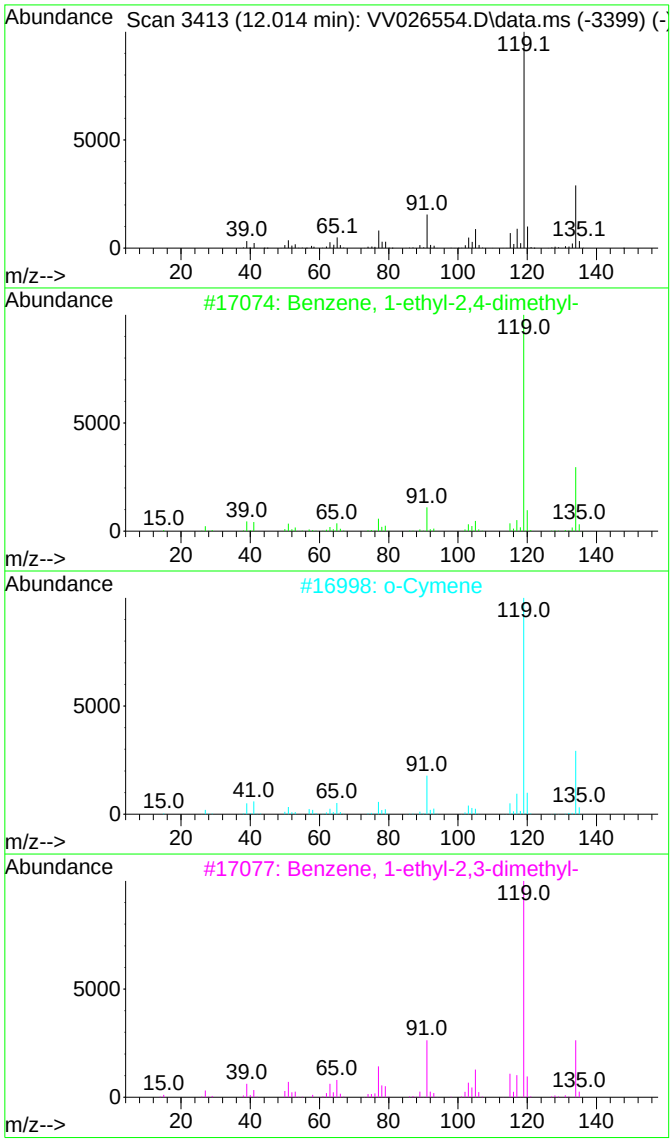
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

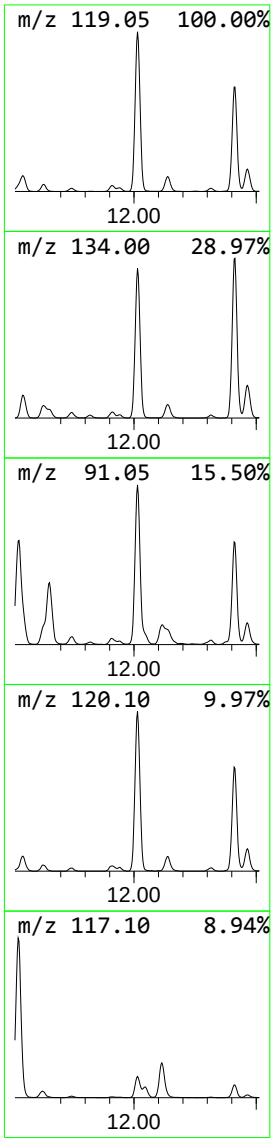
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 17 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.014	17.92 ug/L	2293110	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	97
2	o-Cymene	134	C10H14	000527-84-4	97
3	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	96
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

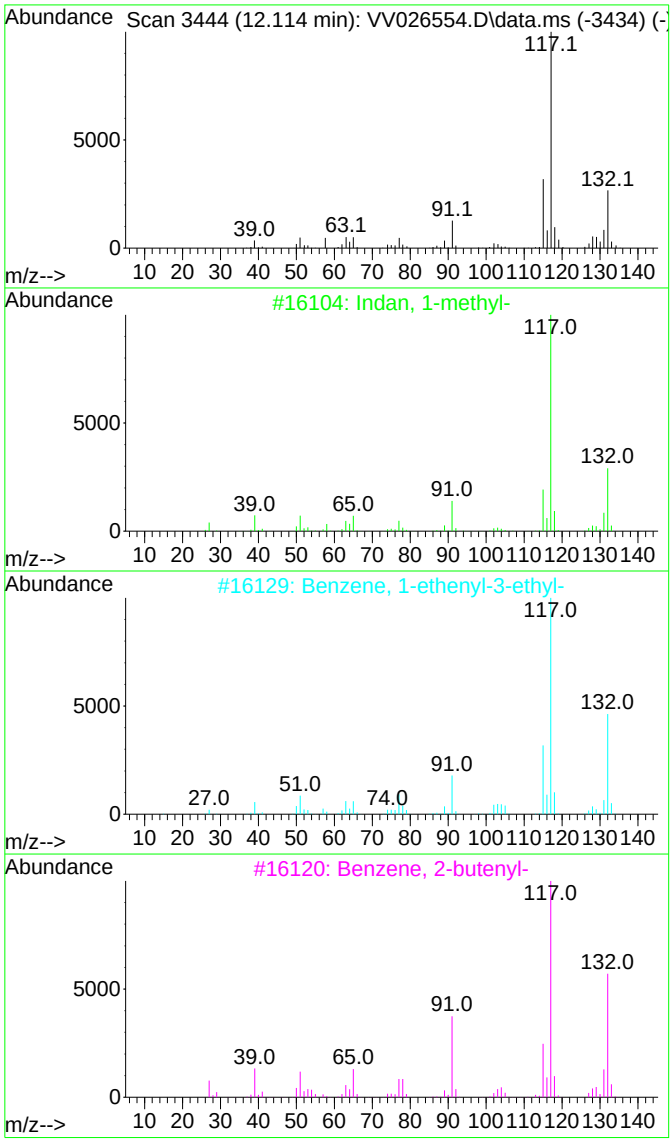
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

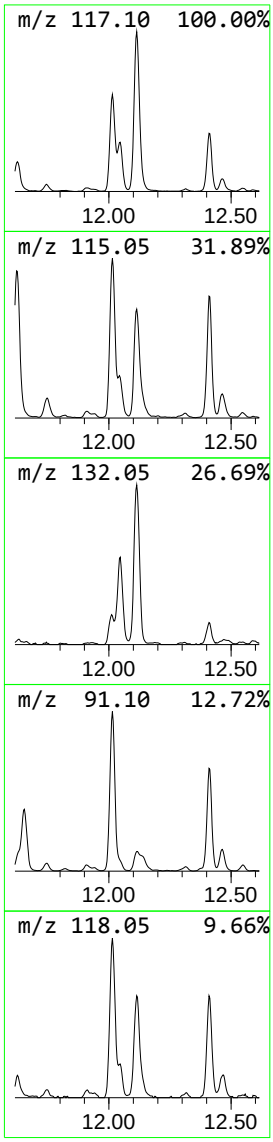
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 18 Indan, 1-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.114	4.66 ug/L	596186	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	87
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	86
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
4	1-Phenyl-1-butene	132	C10H12	000824-90-8	80
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	80





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

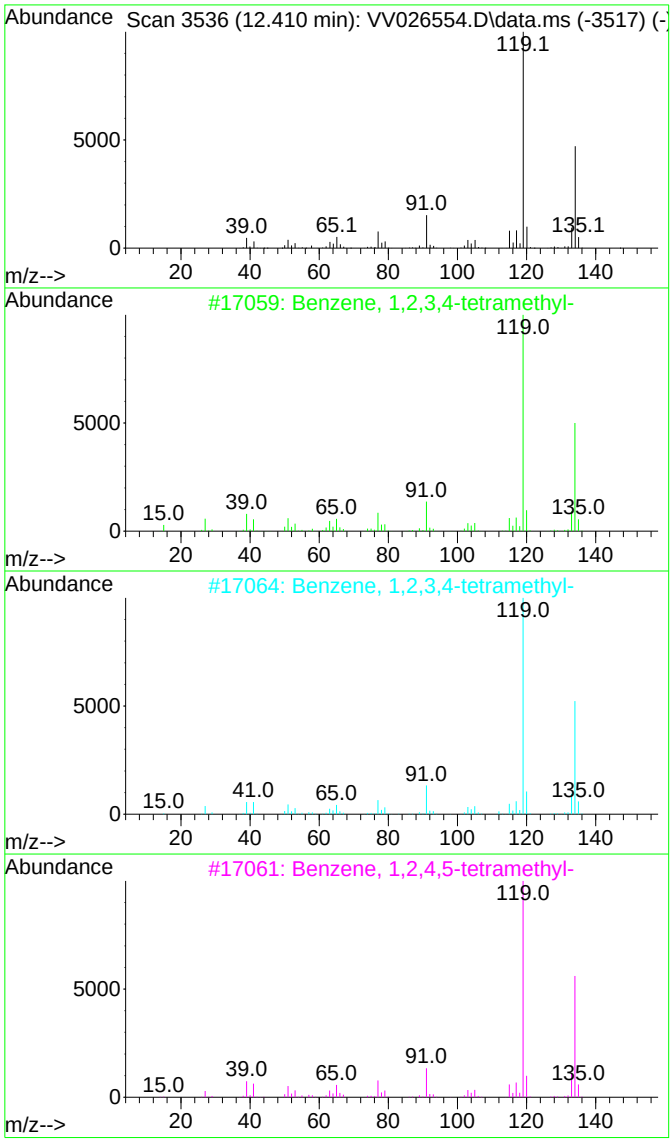
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

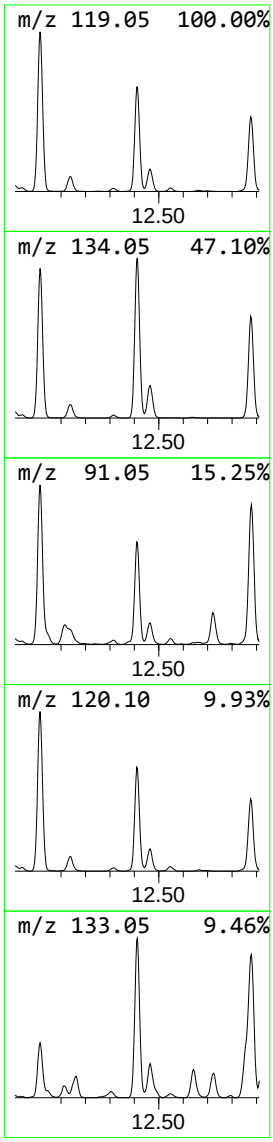
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 19 Benzene, 1,2,3,4-tetramethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.410	12.05 ug/L	1542360	1,4-Dichlorobenzene-d4	11.249

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

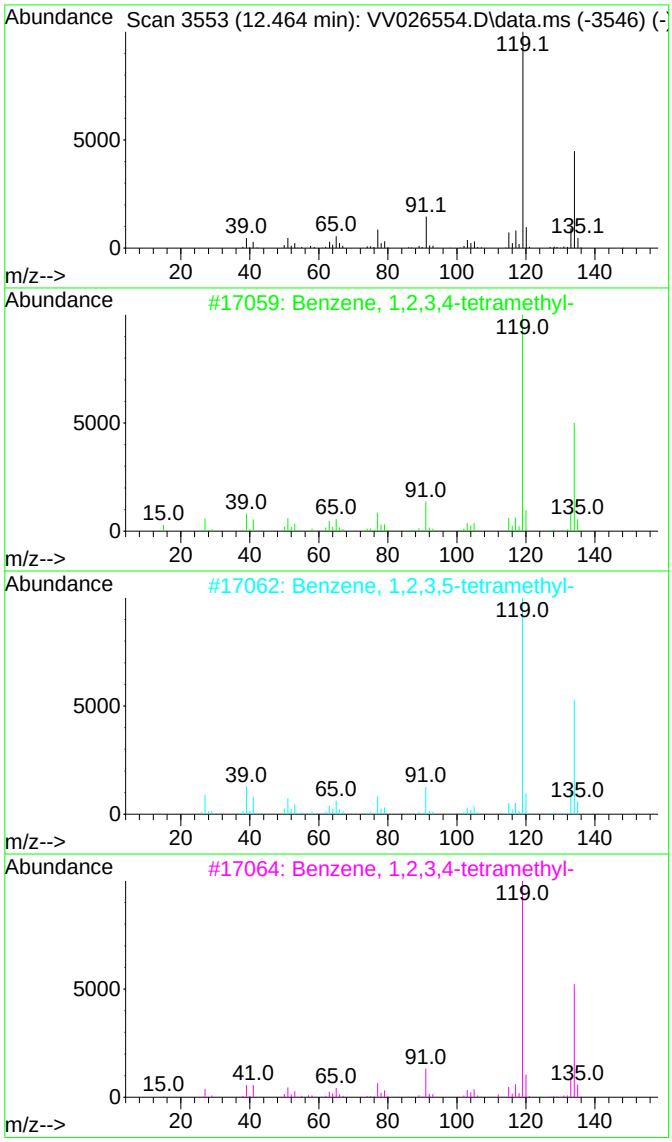
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

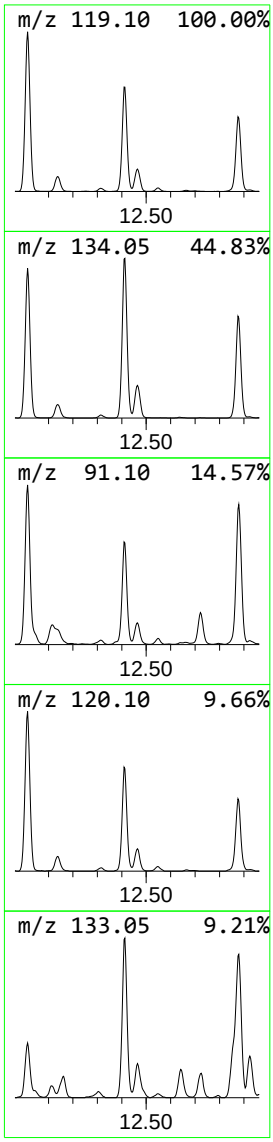
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 20 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	2.74 ug/L	350223	1,4-Dichlorobenzene-d4	11.249

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

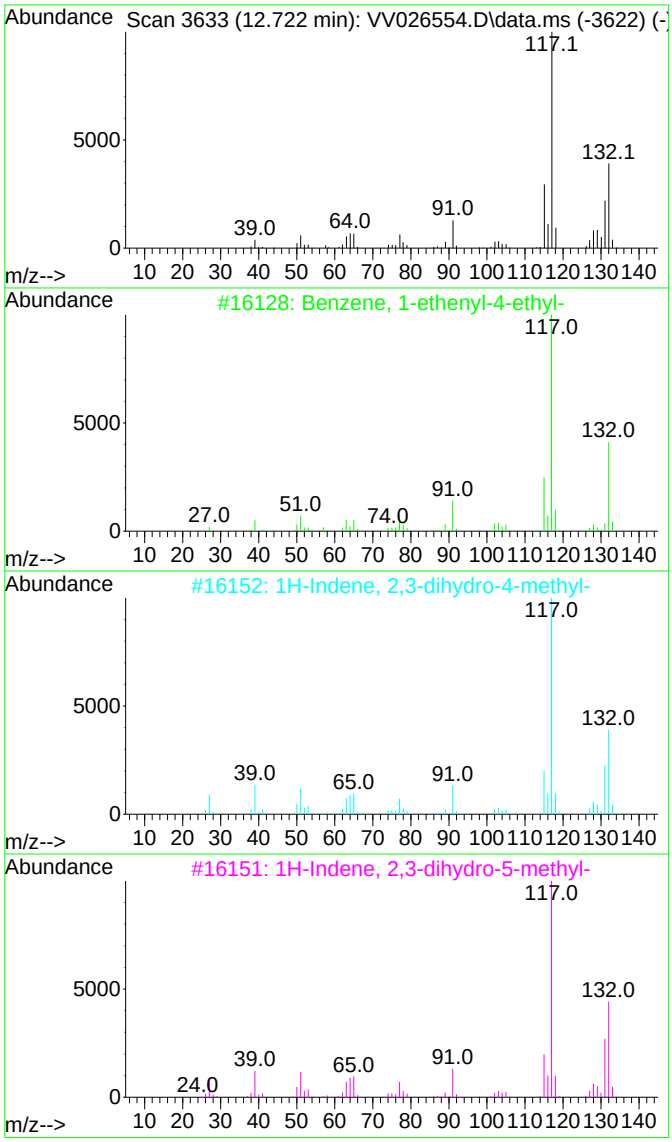
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

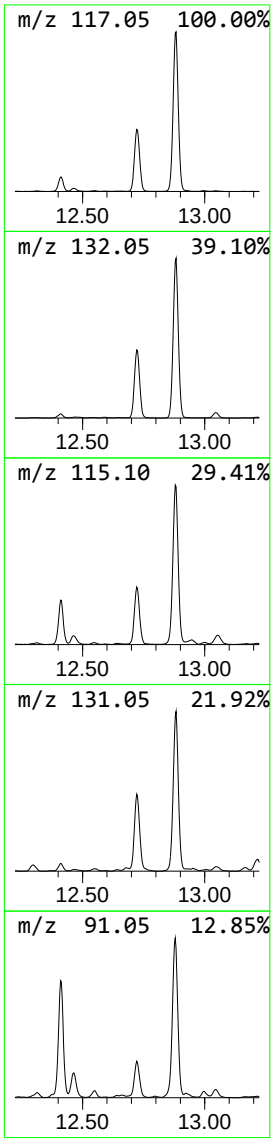
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 21 Benzene, 1-ethenyl-4-ethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.722	5.36 ug/L	686329	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	91
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

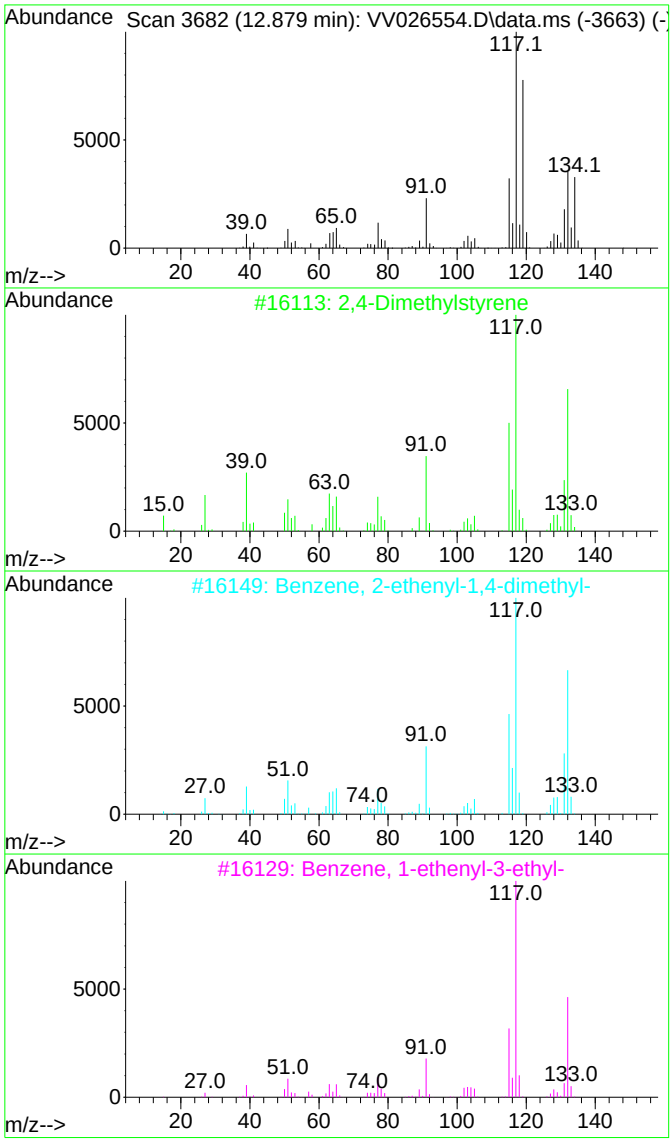
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

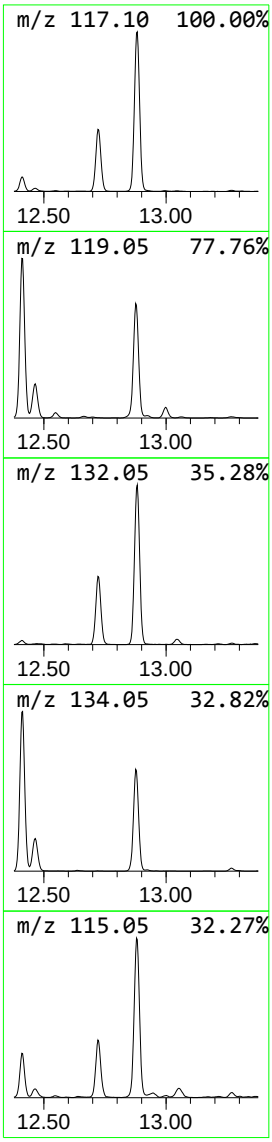
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 22 2,4-Dimethylstyrene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.879	21.03 ug/L	2691050	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
2	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
3	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55
4	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	55
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	55





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

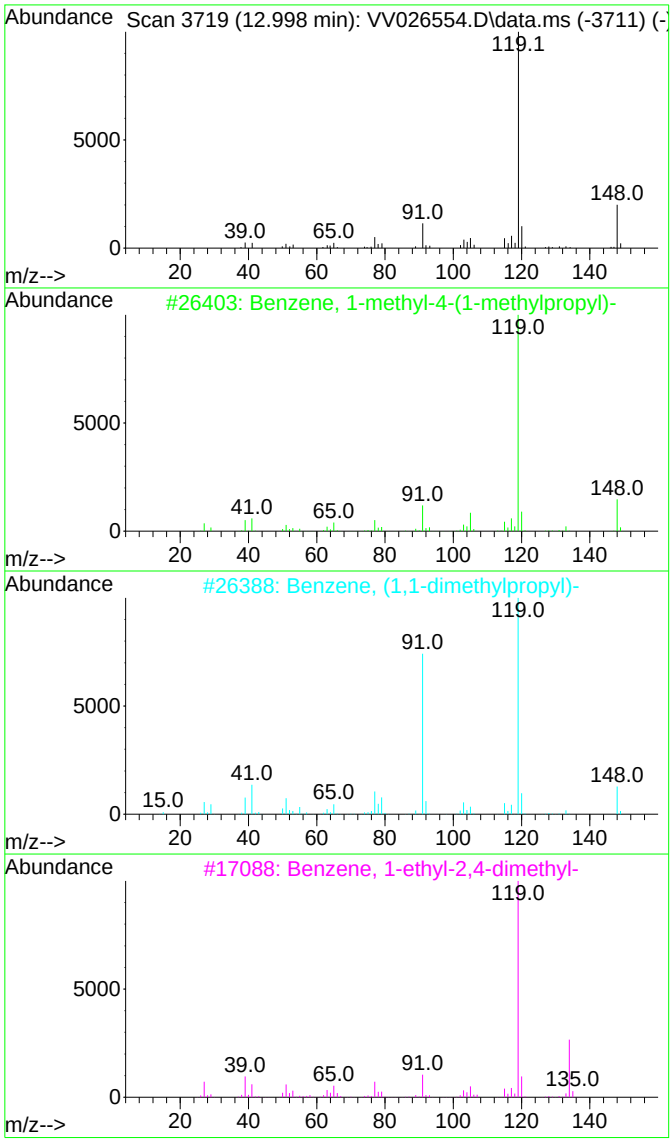
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

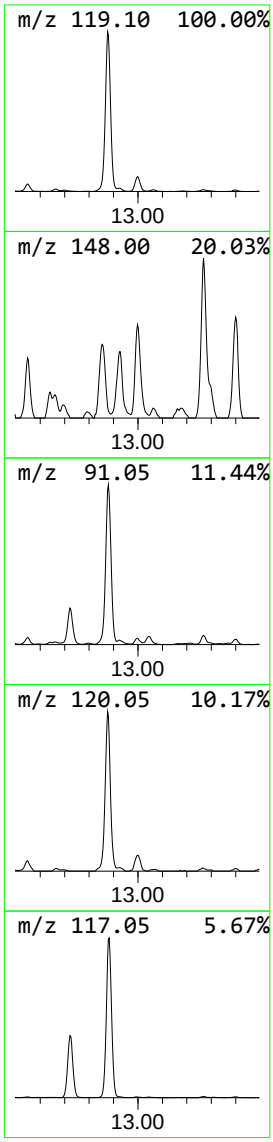
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 23 Benzene, 1-methyl-4-(1-meth... Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.998	0.60 ug/L	76548	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	91
2	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	83
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	83
4	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	78
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

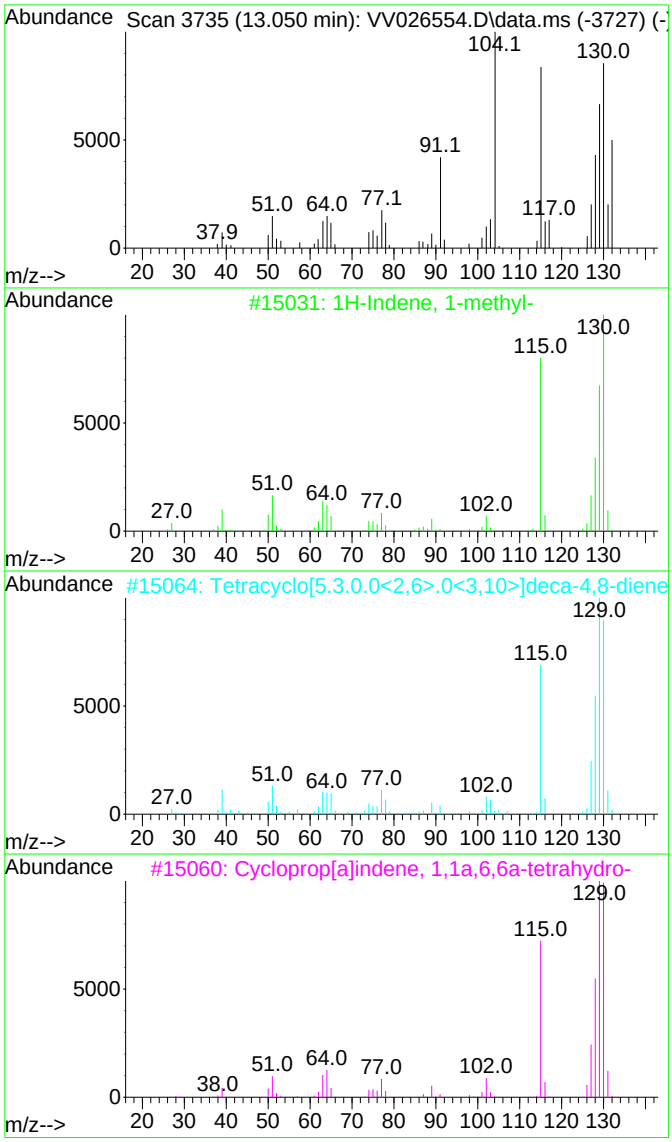
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

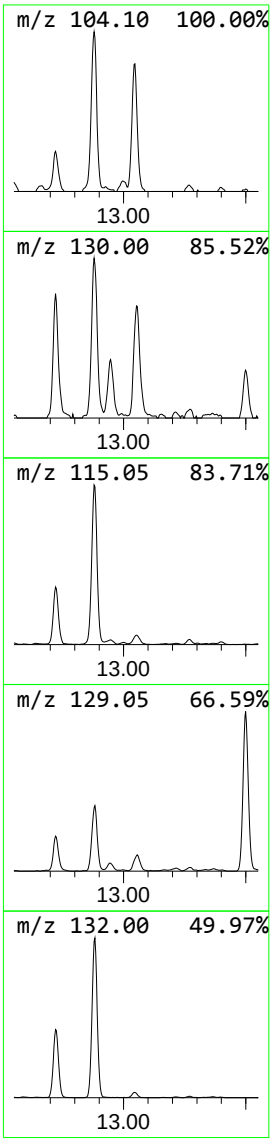
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 24 1H-Indene, 1-methyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.050	0.93 ug/L	119037	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 1-methyl-	130	C10H10	000767-59-9	91
2	Tetracyclo[5.3.0.0<2,6>.0<3,10>]...	130	C10H10	034324-40-8	90
3	Cycloprop[a]indene, 1,1a,6,6a-te...	130	C10H10	015677-15-3	89
4	Benzene, (1-methyl-2-cyclopropen...	130	C10H10	065051-83-4	87
5	2-Methylindene	130	C10H10	002177-47-1	86





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

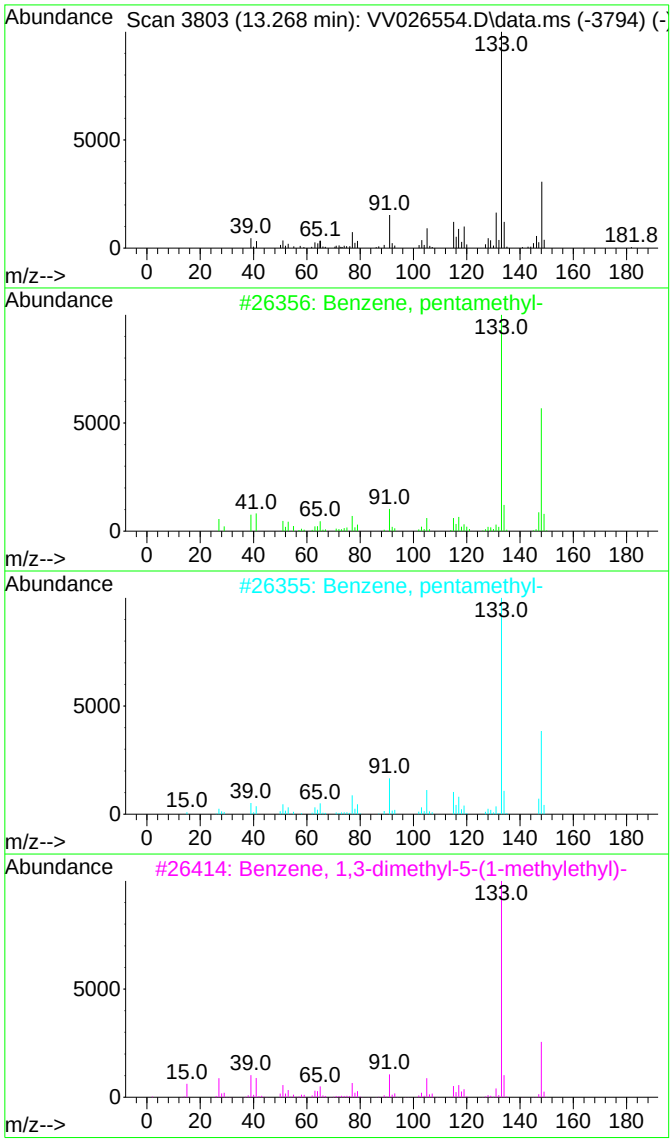
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

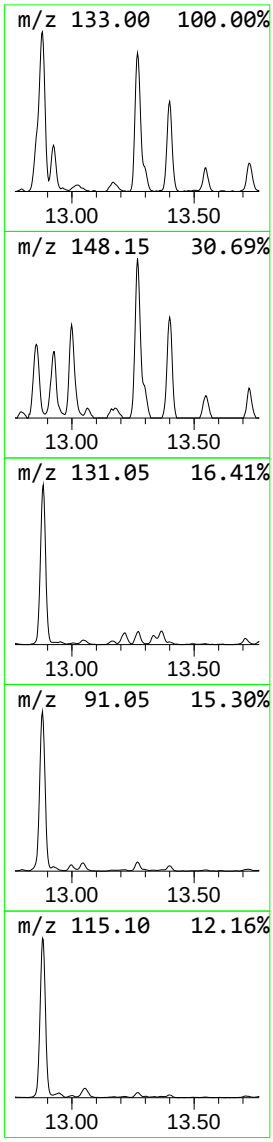
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 25 Benzene, pentamethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.268	1.21 ug/L	154477	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, pentamethyl-	148	C11H16	000700-12-9	87
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	81
3	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	81
4	Benzene, 1-(1,1-dimethylethyl)-3...	148	C11H16	001075-38-3	76
5	4-tert-Butyltoluene	148	C11H16	000098-51-1	76





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

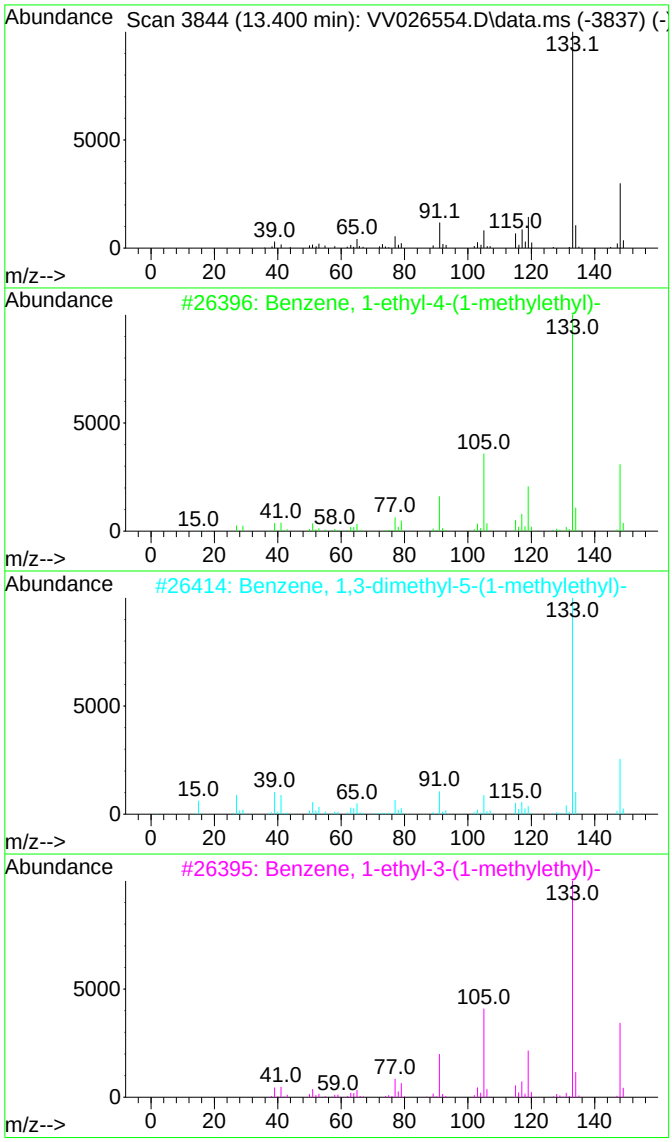
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

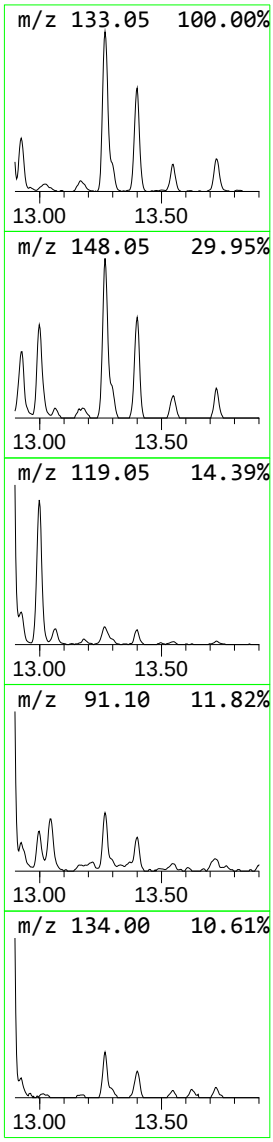
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 26 Benzene, 1-ethyl-4-(1-methy... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.400	0.64 ug/L	81297	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	94
2	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	91
3	Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	90
4	Benzene, 1-ethyl-2,4,5-trimethyl-	148	C11H16	017851-27-3	90
5	Benzene, pentamethyl-	148	C11H16	000700-12-9	90





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

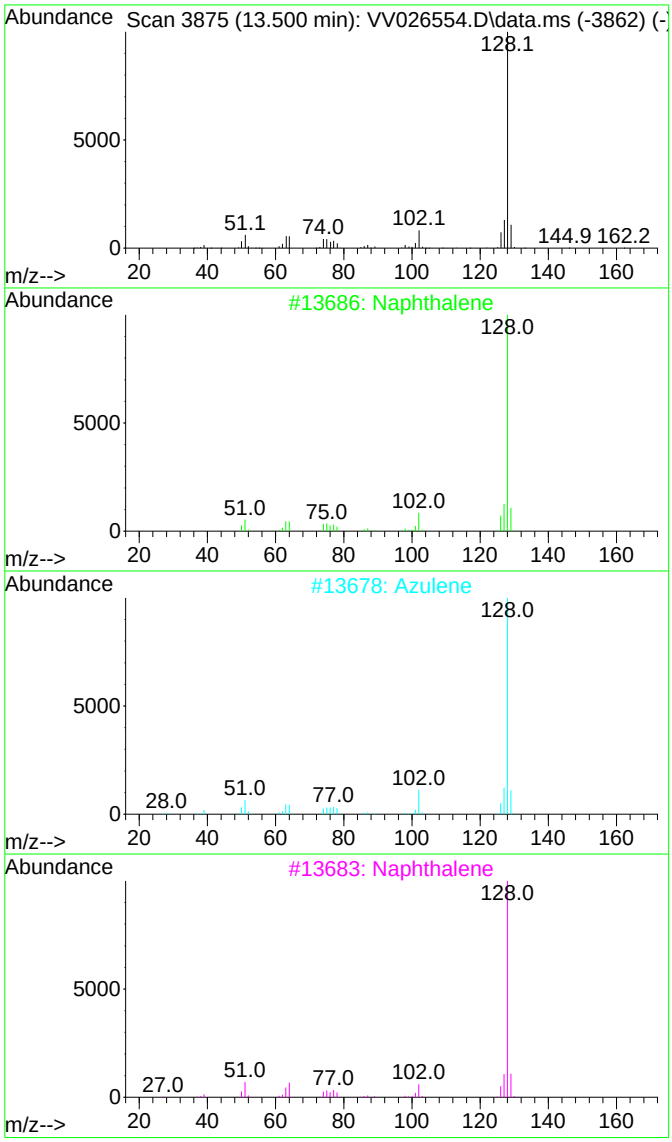
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

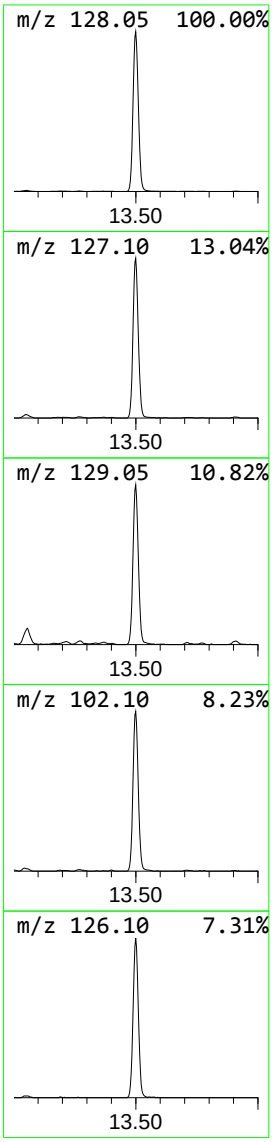
****

Peak Number 27 Azulene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.500	11.64 ug/L	1489810	1,4-Dichlorobenzene-d4	11.249

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





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

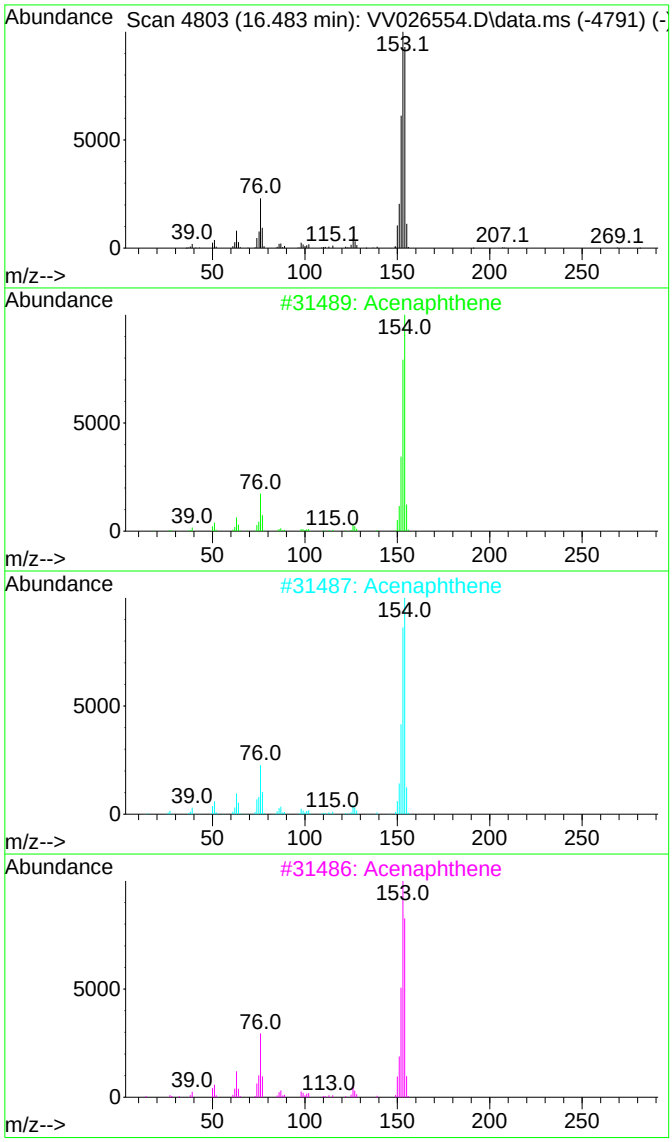
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

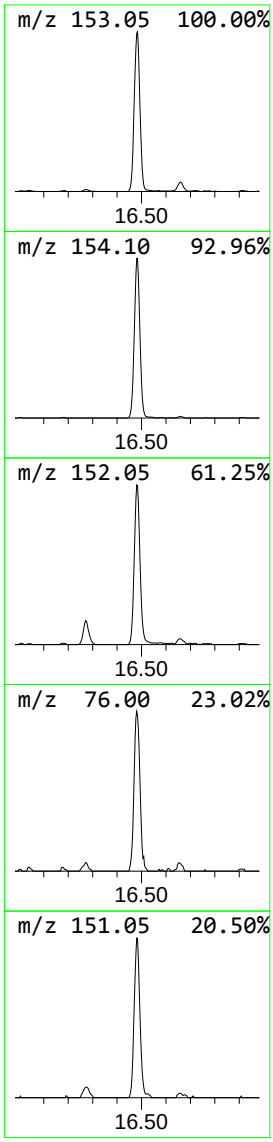
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 28 Hexa-2,4-diyn-1-ylbenzene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.483	2.63 ug/L	335968	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Acenaphthene	154	C12H10	000083-32-9	93
2	Acenaphthene	154	C12H10	000083-32-9	90
3	Acenaphthene	154	C12H10	000083-32-9	86
4	Acenaphthene	154	C12H10	000083-32-9	74
5	Hexa-2,4-diyn-1-ylbenzene	154	C12H10	000520-74-1	53





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV062322\
Data File : VV026554.D
Acq On : 24 Jun 2022 14:39
Operator : SY/MD
Sample : N3401-14DL 10X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0AF2DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR062322WMA.M
Quant Title : TRACE VOA SFAM1.0

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	1.449	2.6	ug/L	305652	1	5.625	585509	5.0
(DEL) Alkane: C...	3.777	3.0	ug/L	356208	1	5.625	585509	5.0
(DEL) Alkane: S...	4.777	0.9	ug/L	103113	1	5.625	585509	5.0
(DEL) Alkane: C...	4.957	1.1	ug/L	131440	1	5.625	585509	5.0
(DEL) Alkane: C...	5.185	1.7	ug/L	197830	1	5.625	585509	5.0
(DEL) Alkane: C...	5.259	1.6	ug/L	186742	1	5.625	585509	5.0
(DEL) Alkane: C...	5.333	2.0	ug/L	235689	1	5.625	585509	5.0
(DEL) Alkane: C...	6.368	0.7	ug/L	79260	1	5.625	585509	5.0
(DEL) Alkane: C...	7.268	0.5	ug/L	66547	2	8.854	641207	5.0
Benzene, propyl-	10.352	35.7	ug/L	4571520	3	11.249	639939	5.0
Benzene, (1-met...	11.088	0.8	ug/L	97528	3	11.249	639939	5.0
Benzene, 1,2,3-...	11.336	2.0	ug/L	262755	3	11.249	639939	5.0
Indane	11.525	16.7	ug/L	2140150	3	11.249	639939	5.0
Benzene, 1,2-di...	11.744	0.9	ug/L	116974	3	11.249	639939	5.0
Benzene, 2-ethy...	11.911	0.7	ug/L	92040	3	11.249	639939	5.0
Benzene, 1-ethy...	12.014	17.9	ug/L	2293110	3	11.249	639939	5.0
Indan, 1-methyl-	12.114	4.7	ug/L	596186	3	11.249	639939	5.0
Benzene, 1,2,3,...	12.410	12.1	ug/L	1542360	3	11.249	639939	5.0
Benzene, 1,2,3,...	12.464	2.7	ug/L	350223	3	11.249	639939	5.0
Benzene, 1-ethe...	12.722	5.4	ug/L	686329	3	11.249	639939	5.0
2,4-Dimethylsty...	12.879	21.0	ug/L	2691050	3	11.249	639939	5.0
Benzene, 1-meth...	12.998	0.6	ug/L	76548	3	11.249	639939	5.0
1H-Indene, 1-me...	13.050	0.9	ug/L	119037	3	11.249	639939	5.0
Benzene, pentam...	13.268	1.2	ug/L	154477	3	11.249	639939	5.0
Benzene, 1-ethy...	13.400	0.6	ug/L	81297	3	11.249	639939	5.0
Azulene	13.500	11.6	ug/L	1489810	3	11.249	639939	5.0
Hexa-2,4-diyn-1...	16.483	2.6	ug/L	335968	3	11.249	639939	5.0