LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

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_U\Method\SFAMUTR102822WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU051647.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.375	311	322	342	rBV	3825721	5702930	6.26%	1.779%
2	3.356	612	627	649	rVB	437271	967665	1.06%	0.302%
3	3.694	716	732	751	rVB	424041	930933	1.02%	0.290%
4	3.990	805	824	855	rBV	4955629	12678554	13.92%	3.956%
5	4.526	974	991	1013	rBV	1528876	3719467	4.08%	1.160%
6	5.063	1140	1158	1213	rBV2	492213	1411402	1.55%	0.440%
7	5.385	1241	1258	1273	rBV	464688	1069527	1.17%	0.334%
8	5.777	1345	1380	1397	rBV2	41595925	91110990	100.00%	28.426%
9	6.761	1674	1686	1707	rVB	582221	1205620	1.32%	0.376%
10	7.967	2050	2061	2076	rBV	1630277	2716129	2.98%	0.847%
11	8.632	2256	2268	2298	rBV	778759	1542536	1.69%	0.481%
12	9.449	2499	2522	2547	rBV2	39833565	66469388	72.95%	20.738%
13	9.568	2548	2559	2582	rVB2	5271469	9261852	10.17%	2.890%
14	9.690	2583	2597	2612	rBV2	27816253	44156330	48.46%	13.776%
15	10.099	2710	2724	2749	rVB	9407458	14927343	16.38%	4.657%
16	10.481	2829	2843	2856	rBV	4518459	7021964	7.71%	2.191%
17	10.902	2956	2974	2988	rBV	3096710	5812518	6.38%	1.813%
18	10.989	2989	3001	3009	rBV2	2134235	4078167	4.48%	1.272%
19	11.086	3022	3031	3047	rVB2	1156950	1879778	2.06%	0.586%
20	11.288	3083	3094	3117	rVB	2708422	4286879	4.71%	1.337%
21	11.465	3139	3149	3162	rVB	4321669	6541881	7.18%	2.041%
22	11.832	3243	3263	3272	rVV3	1126412	2821399	3.10%	0.880%
23	11.893	3272	3282	3293	rVB	3436066	5393176	5.92%	1.683%
24	12.095	3326	3345	3362	rVV2	4645506	7603022	8.34%	2.372%
25	12.185	3362	3373	3396	rVB6	1269546	3288489	3.61%	1.026%
26	12.475	3449	3463	3483	rBV3	1022375	2571900	2.82%	0.802%
27	12.568	3483	3492	3501	rBV	1246815	1766321	1.94%	0.551%
28	12.819	3563	3570	3575	rVV	991024	1500138	1.65%	0.468%
29	12.857	3575	3582	3601	rVB2	2077351	3325972	3.65%	1.038%
30	12.970	3603	3617	3626	rBV	840389	1345779	1.48%	0.420%
31	13.449	3756	3766	3781	rVB3	1049644	1860274	2.04%	0.580%
32	16.886	4823	4835	4847	rBV	1036355	1556209	1.71%	0.486%

Sum of corrected areas: 320524532

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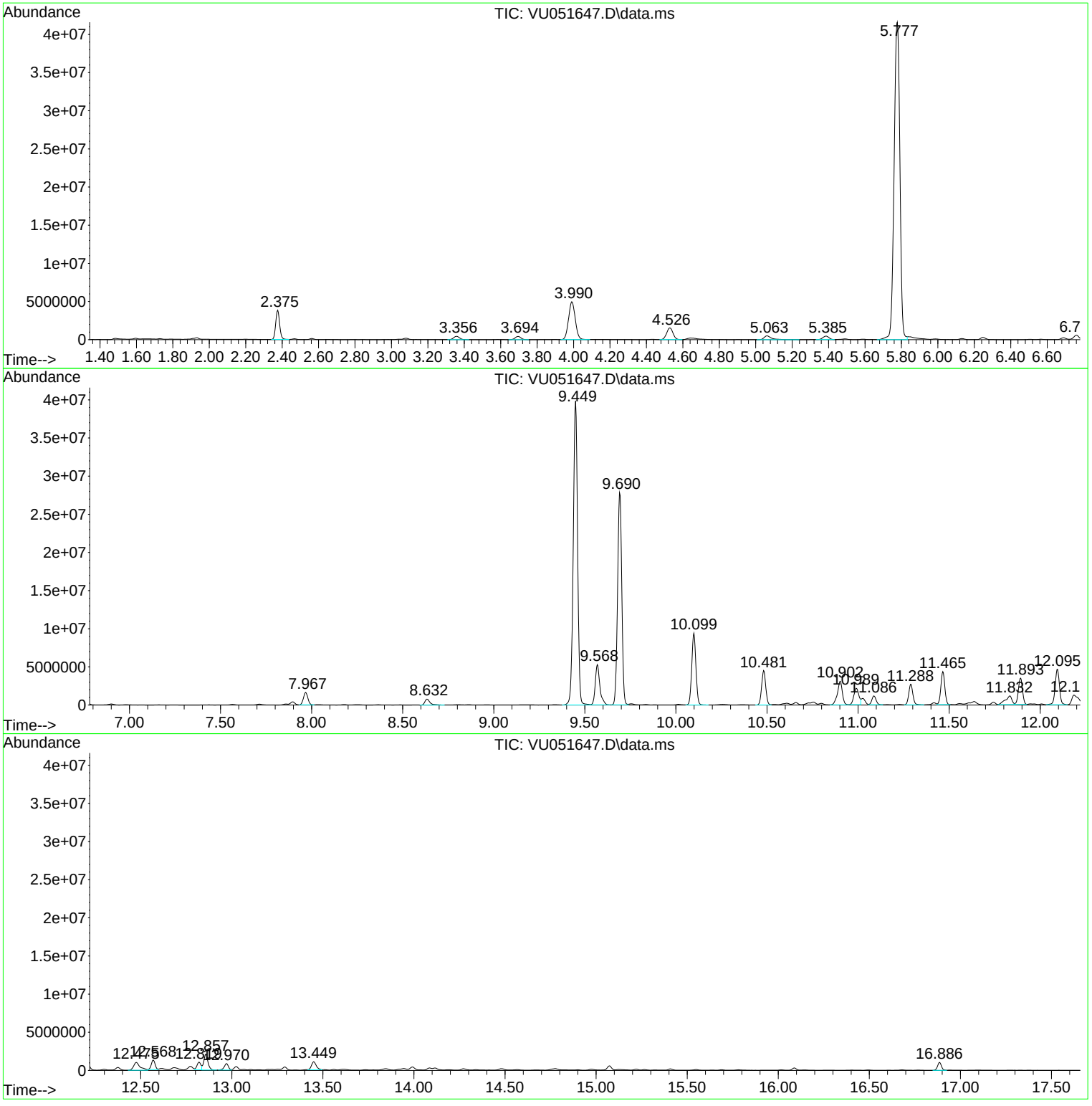
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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Data File : VU051647.D
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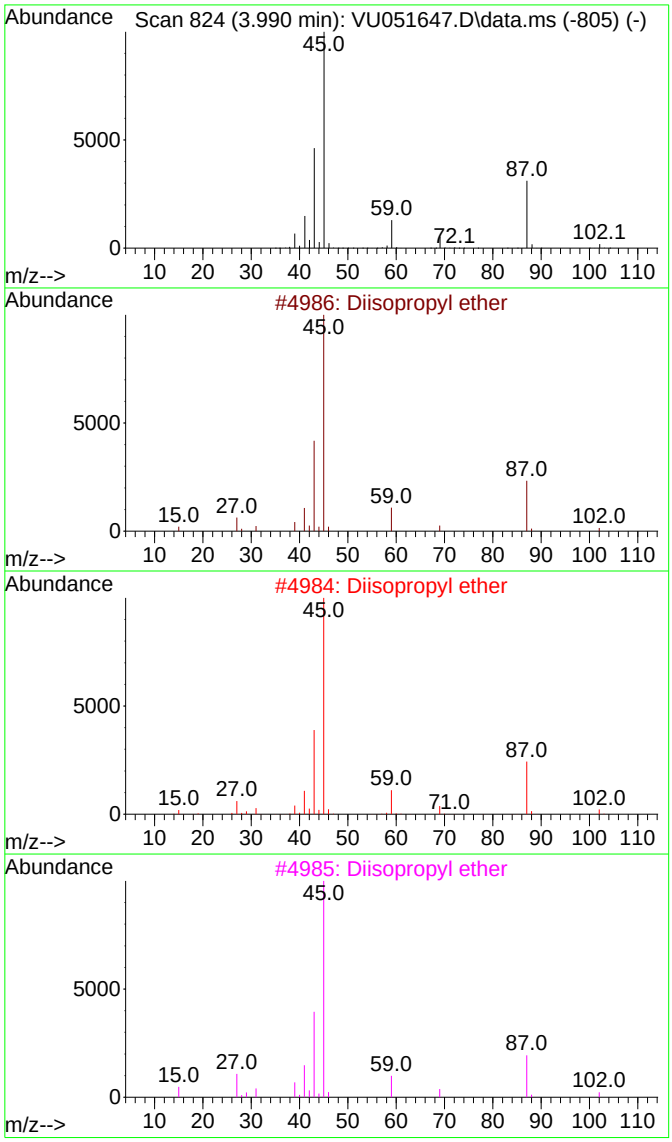
Quant Title : TRACE VOA SFAM1.0

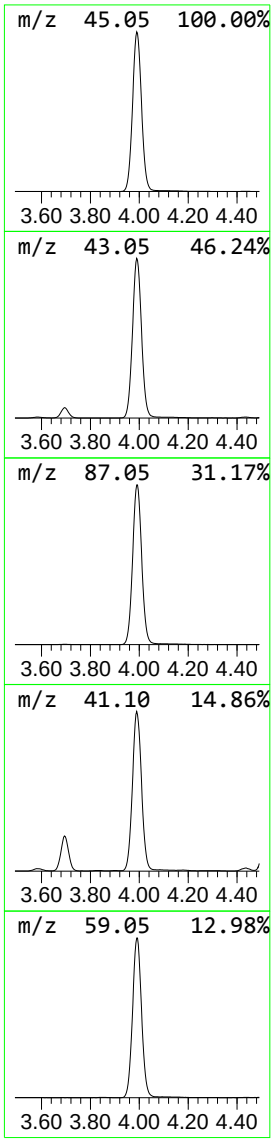
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 Diisopropyl ether Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.990	0.70 ug/L	12678600	1,4-Difluorobenzene	6.247

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Diisopropyl ether	102	C6H14O	000108-20-3	91
2	Diisopropyl ether	102	C6H14O	000108-20-3	91
3	Diisopropyl ether	102	C6H14O	000108-20-3	86
4	Diisopropyl ether	102	C6H14O	000108-20-3	83
5	1-Propanol, 2-(1-methylethoxy)-	118	C6H14O2	003944-37-4	83





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

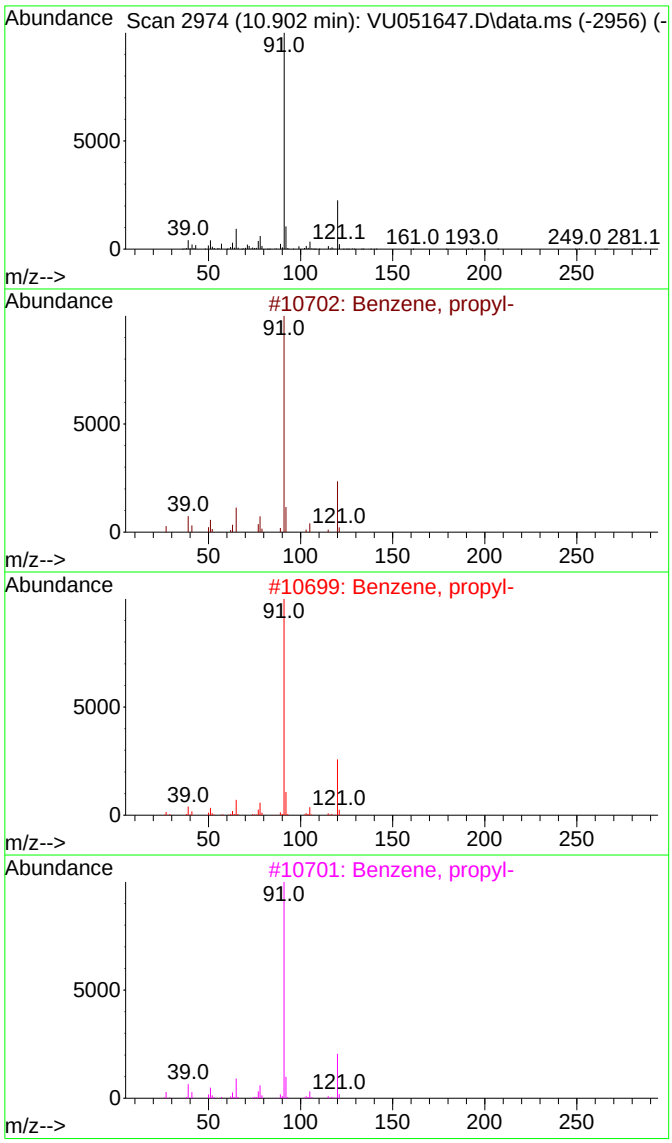
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

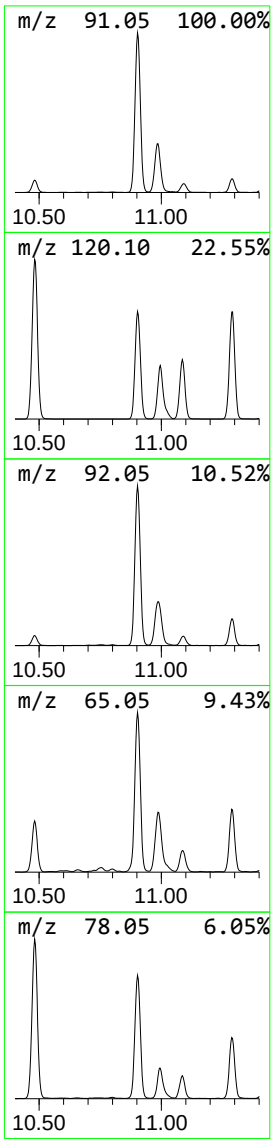
****

Peak Number 2 Benzene, propyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.902	10.30 ug/L	5812520	1,4-Dichlorobenzene-d4	11.809

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-Butylbenzylamine	163	C11H17N	002403-22-7	64
5	Propanenitrile, 3-[(phenylmethyl...	160	C10H12N2	000706-03-6	64





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Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

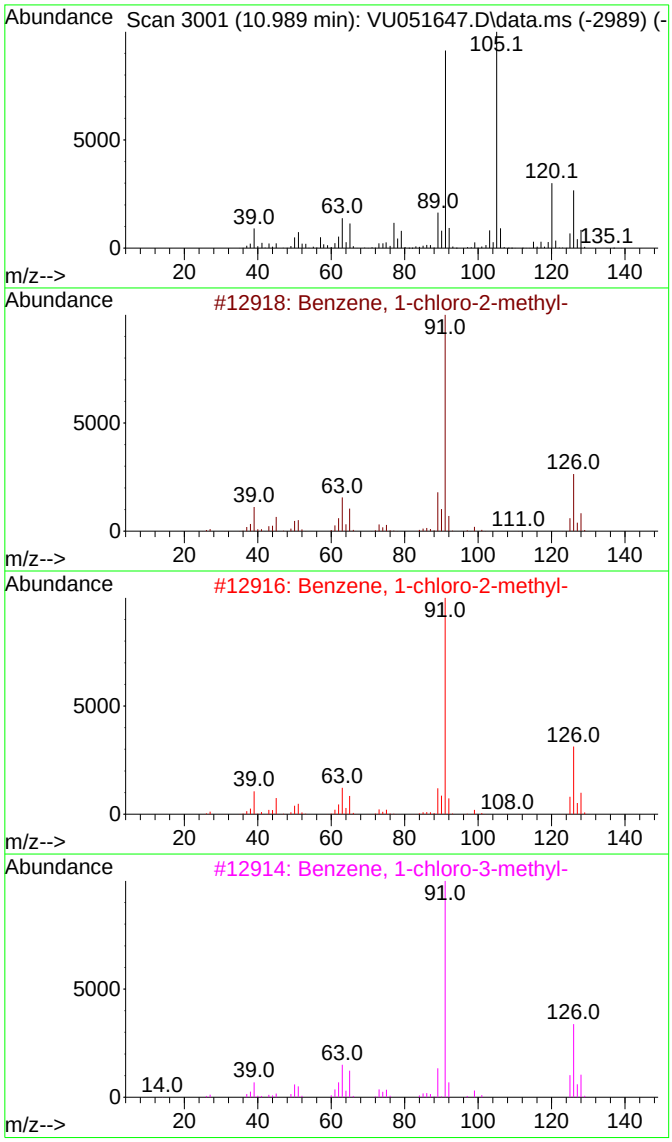
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

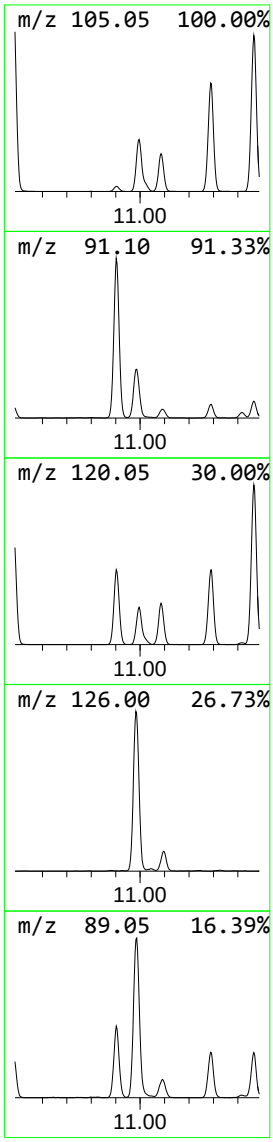
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 Benzene, 1-chloro-2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.989	7.23 ug/L	4078170	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	95
2	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	91
3	Benzene, 1-chloro-3-methyl-	126	C7H7Cl	000108-41-8	91
4	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	90
5	Benzene, 1-chloro-2-methyl-	126	C7H7Cl	000095-49-8	60





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Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

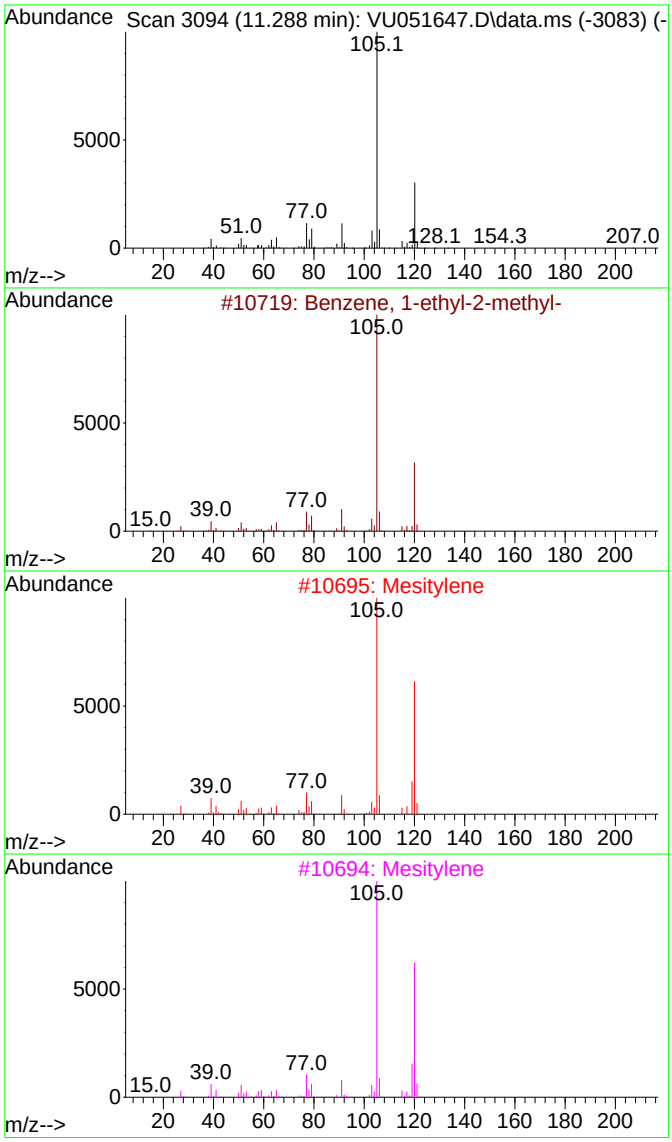
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

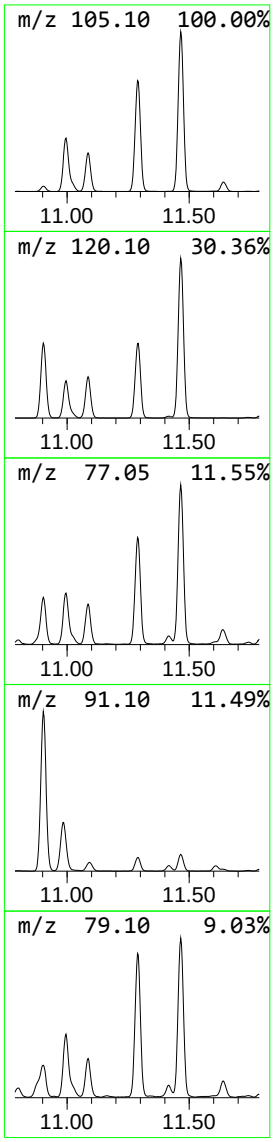
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.288	7.60 ug/L	4286880	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Mesitylene	120	C9H12	000108-67-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

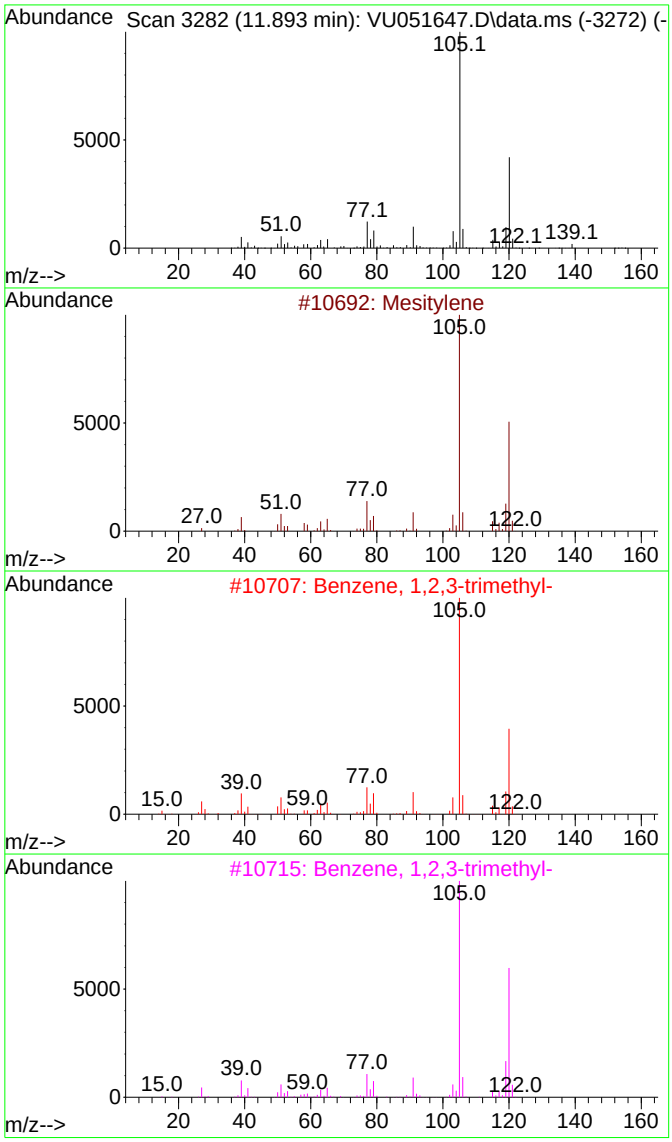
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

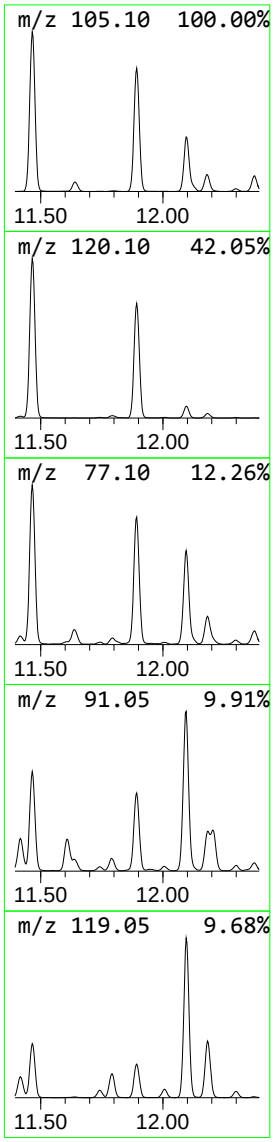
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.893	9.56 ug/L	5393180	1,4-Dichlorobenzene-d4	11.809

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	95
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Mesitylene	120	C9H12	000108-67-8	91





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ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

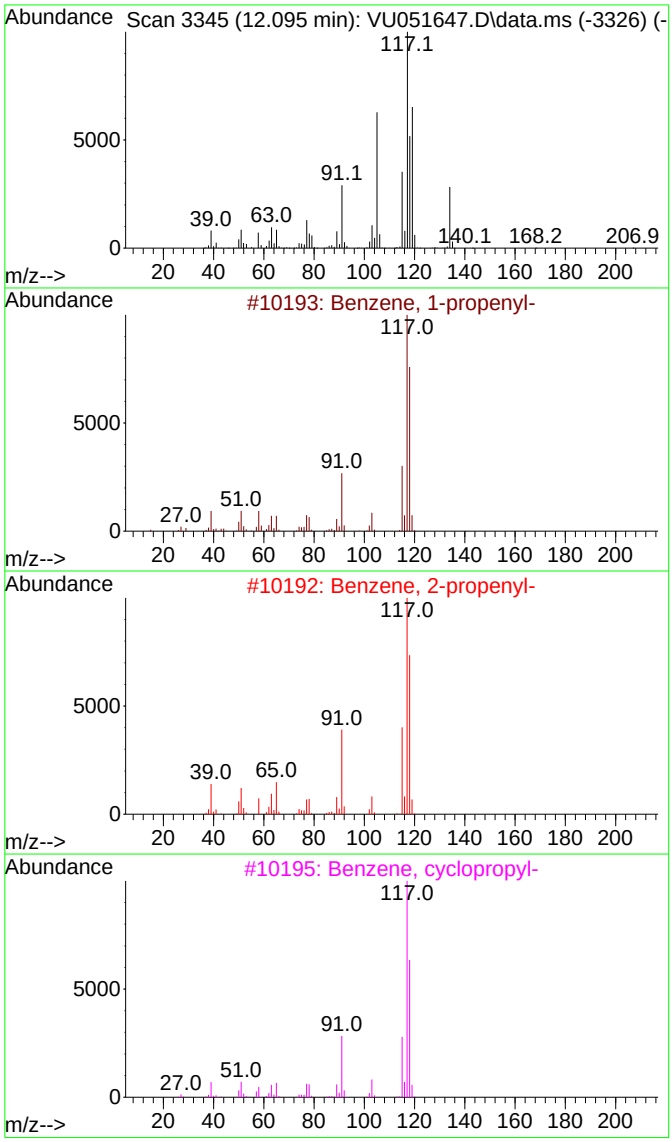
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

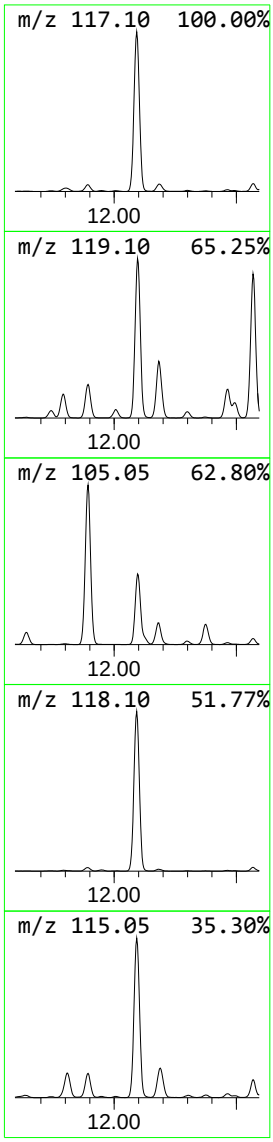
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 Benzene, 1-propenyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	13.47 ug/L	7603020	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
3	Benzene, cyclopropyl-	118	C9H10	000873-49-4	60
4	Benzene, 2-propenyl-	118	C9H10	000300-57-2	60
5	Benzene, 1-propenyl-	118	C9H10	000637-50-3	60





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Operator : JC/MD
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ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

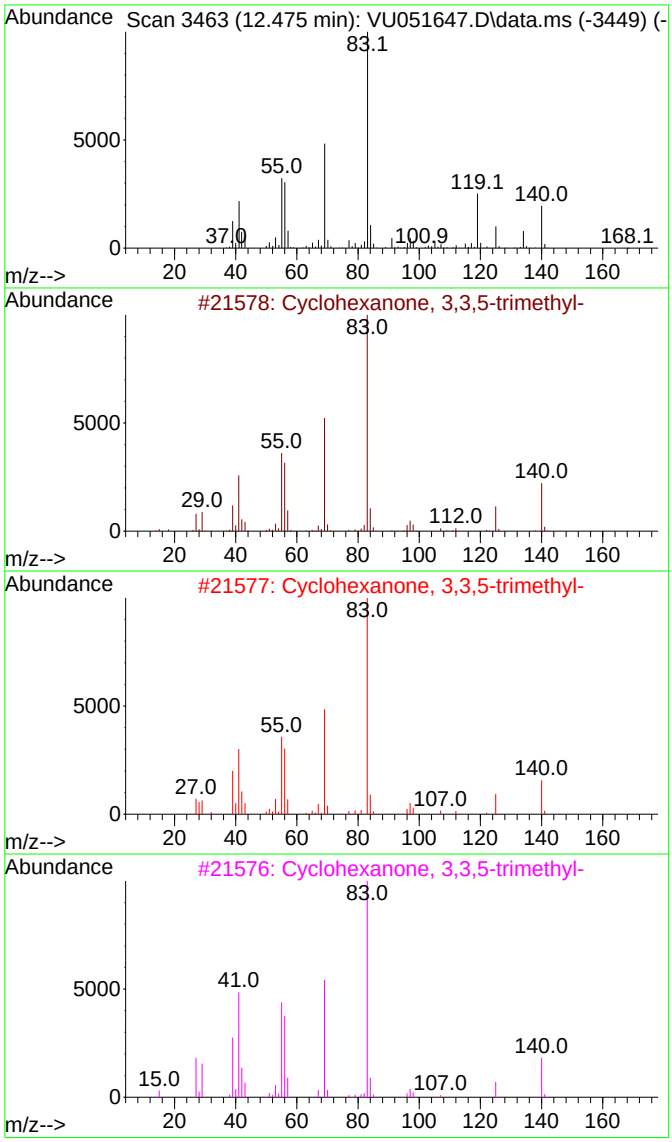
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

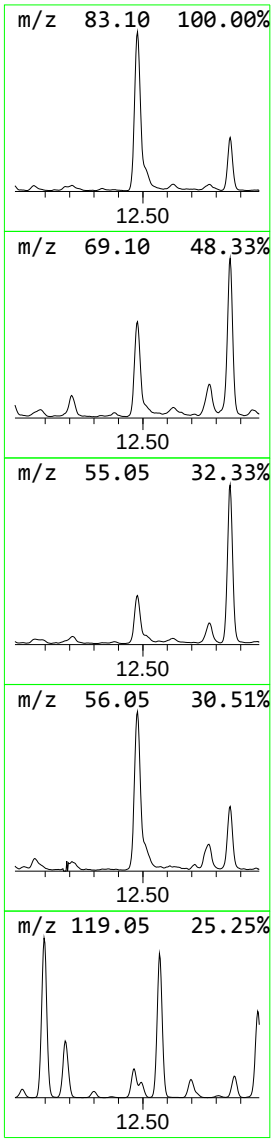
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 Cyclohexanone, 3,3,5-trimet... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.475	4.56 ug/L	2571900	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	96
2	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
3	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4	Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
5	1,3-Cyclohexanedione, 5,5-dimethyl-	140	C8H12O2	000126-81-8	43





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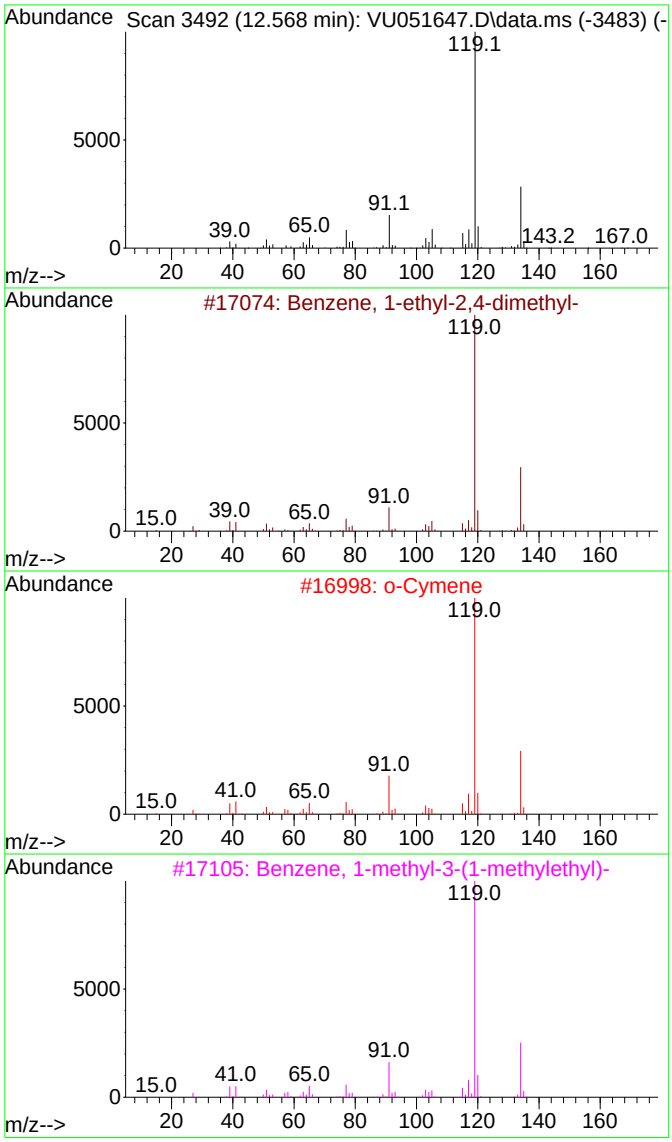
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

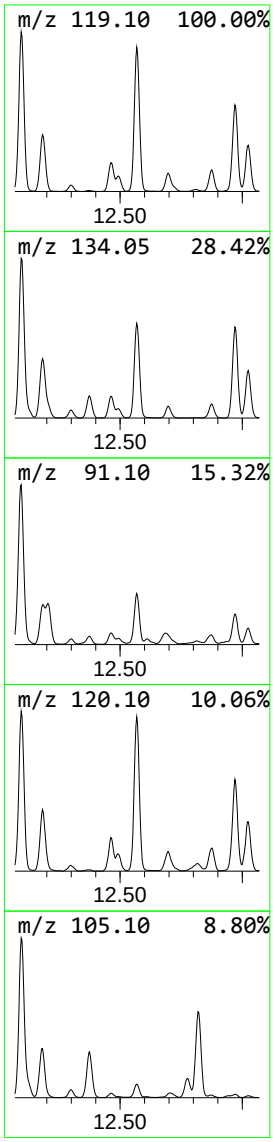
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.568	3.13 ug/L	1766320	1,4-Dichlorobenzene-d4	11.809

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-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
5	o-Cymene	134	C10H14	000527-84-4	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

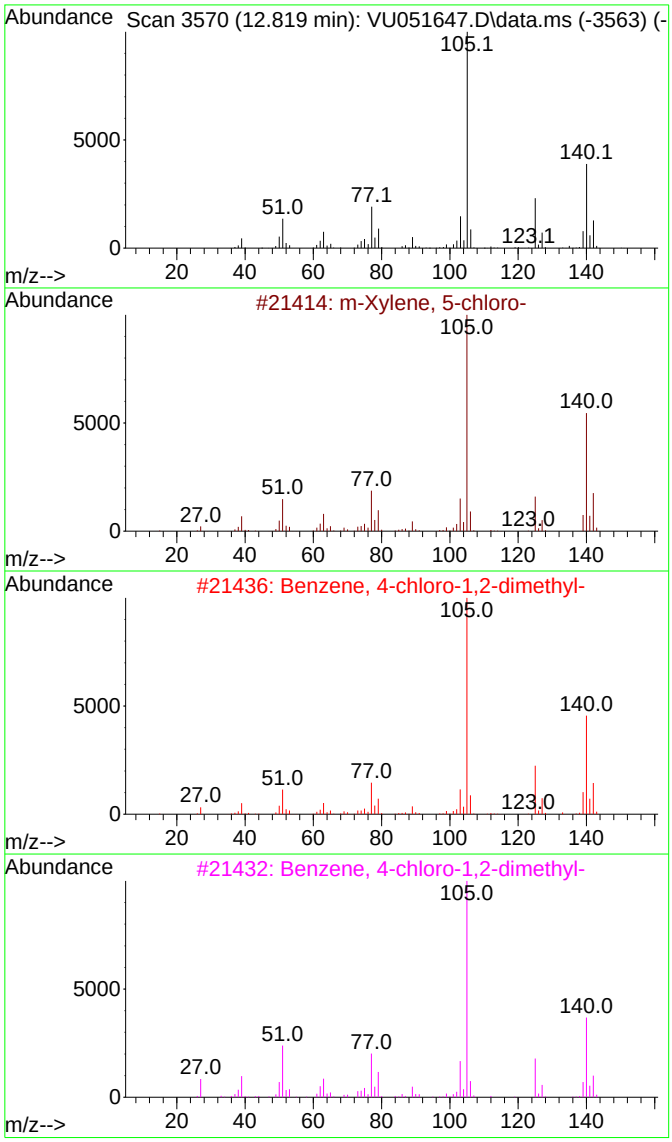
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

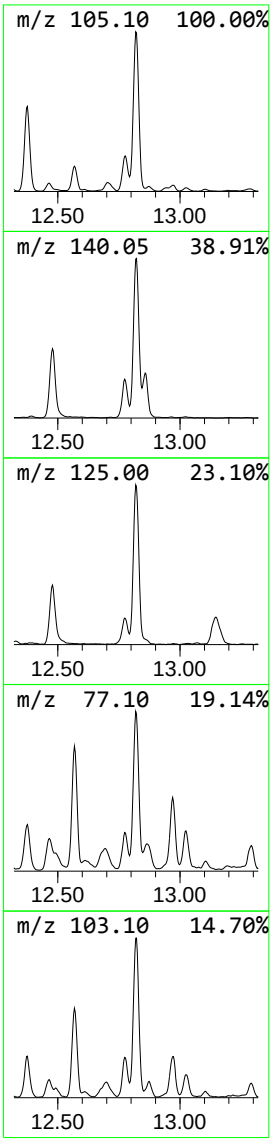
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 m-Xylene, 5-chloro- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.819	2.66 ug/L	1500140	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	m-Xylene, 5-chloro-	140	C8H9Cl	000556-97-8	96
2	Benzene, 4-chloro-1,2-dimethyl-	140	C8H9Cl	000615-60-1	95
3	Benzene, 4-chloro-1,2-dimethyl-	140	C8H9Cl	000615-60-1	94
4	Benzene, 1-chloro-2,3-dimethyl-	140	C8H9Cl	000608-23-1	94
5	Benzene, 4-chloro-1,2-dimethyl-	140	C8H9Cl	000615-60-1	94





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Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

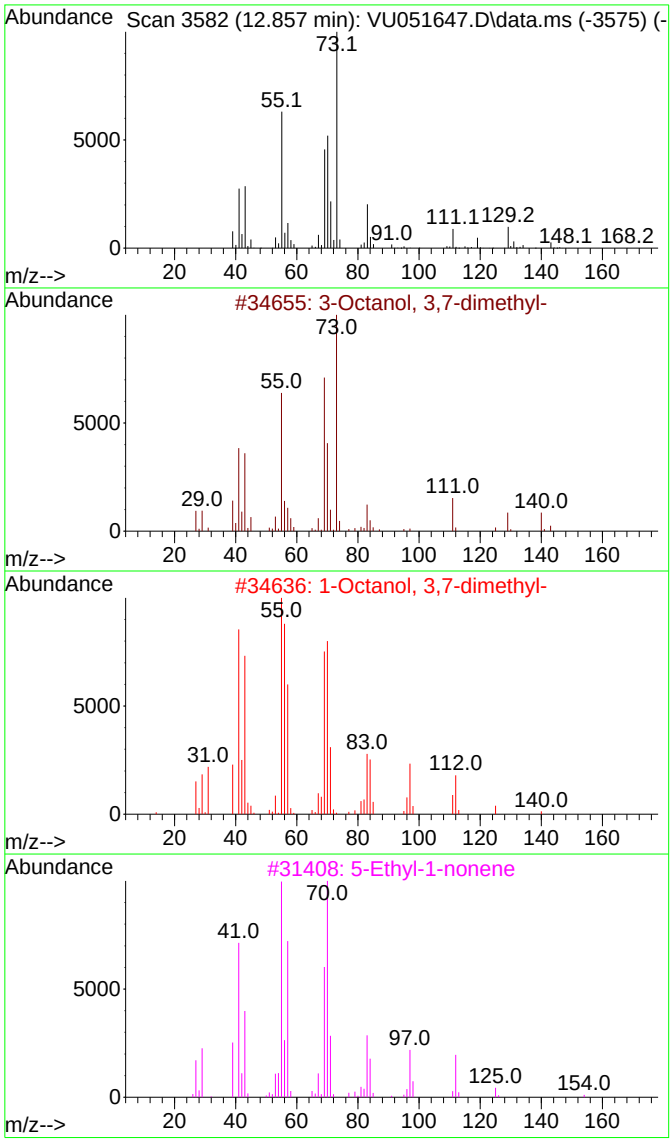
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

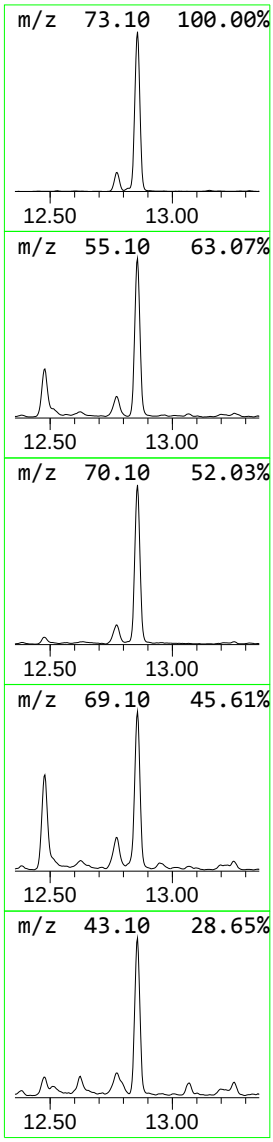
****

Peak Number 10 unknown-01 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.857	5.89 ug/L	3325970	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	40
2	1-Octanol, 3,7-dimethyl-	158	C10H22O	000106-21-8	40
3	5-Ethyl-1-nonene	154	C11H22	019780-74-6	40
4	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	37
5	3-Octanol, 3,7-dimethyl-	158	C10H22O	000078-69-3	35





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

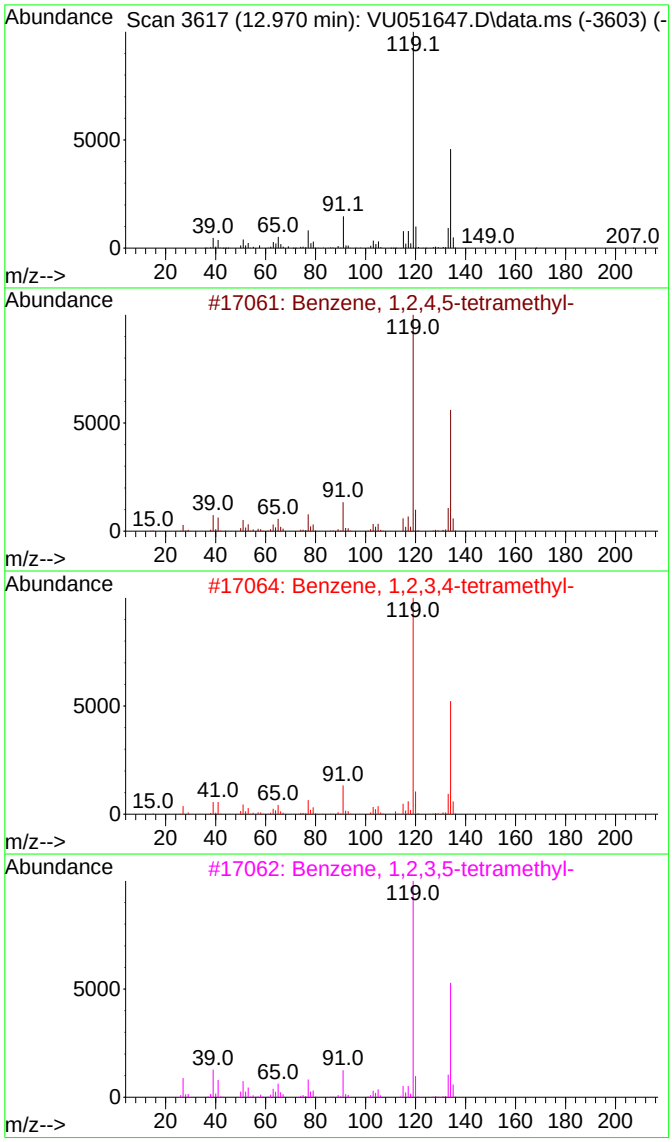
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

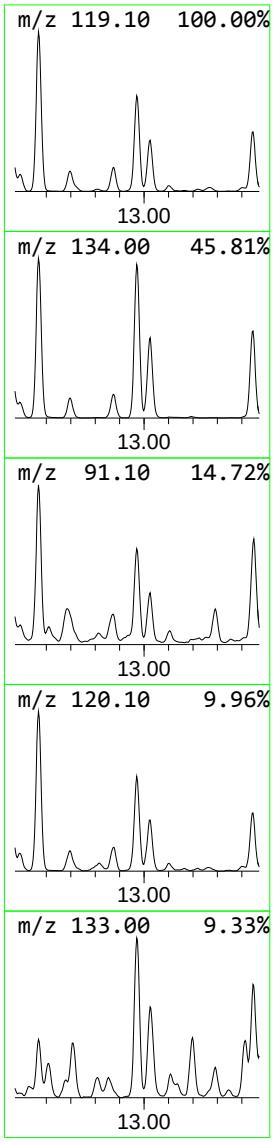
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.970	2.38 ug/L	1345780	1,4-Dichlorobenzene-d4	11.809

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	97
4	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	97
5	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

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

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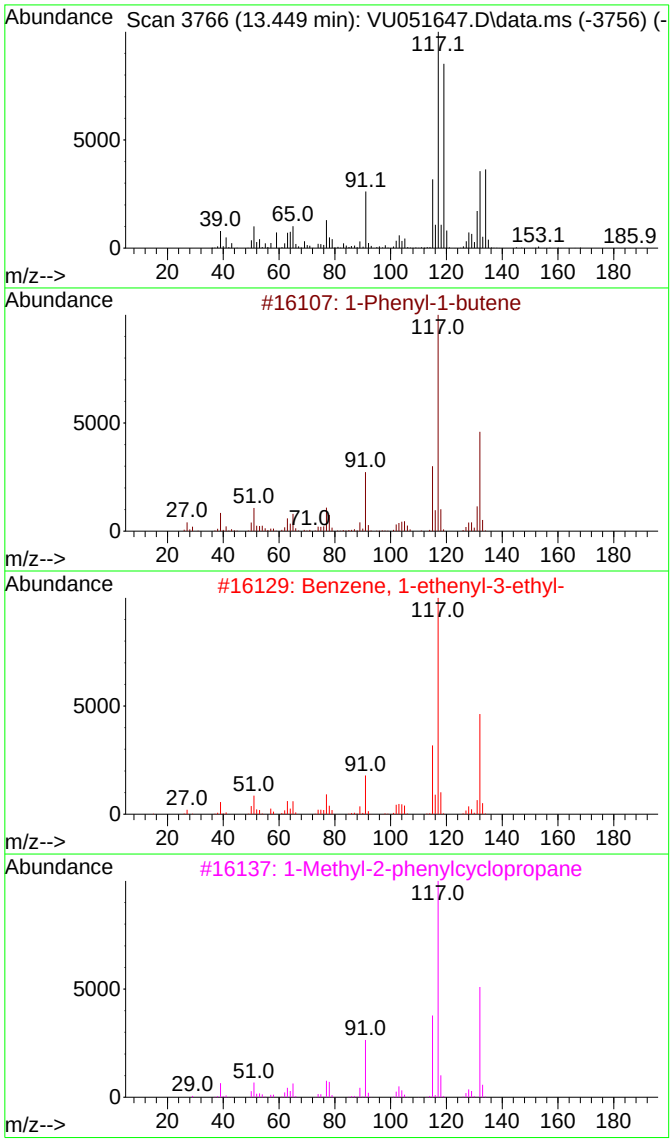
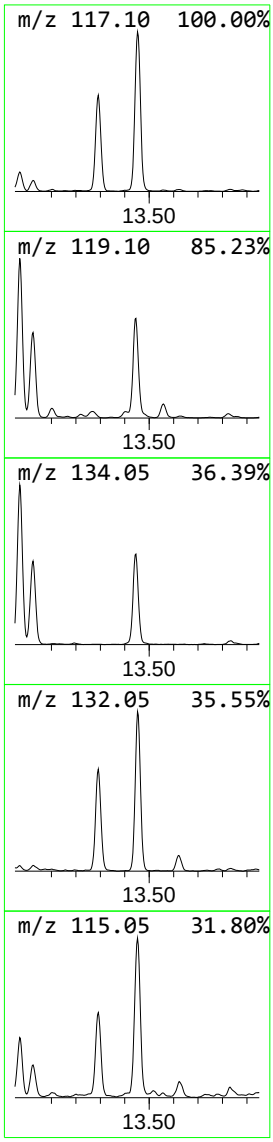
Peak Number 12 1-Phenyl-1-butene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.449	3.30 ug/L	1860270	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	92
2	Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	87
3	1-Methyl-2-phenylcyclopropane	132	C10H12	003145-76-4	86
4	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	86
5	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	62





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

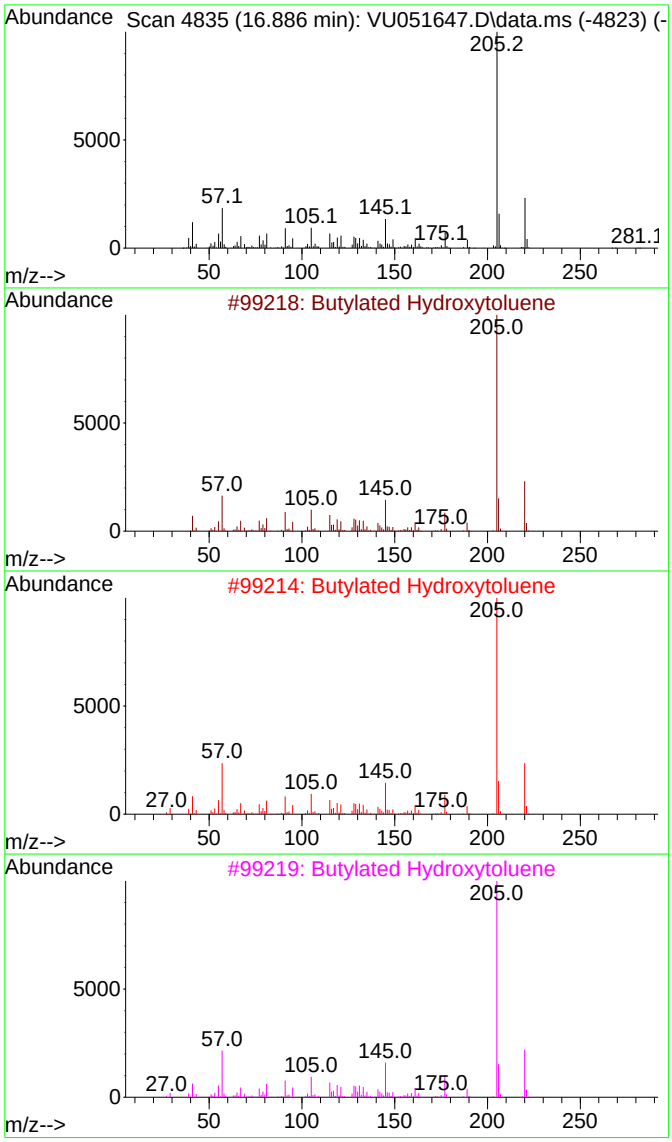
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.M
Quant Title : TRACE VOA SFAM1.0

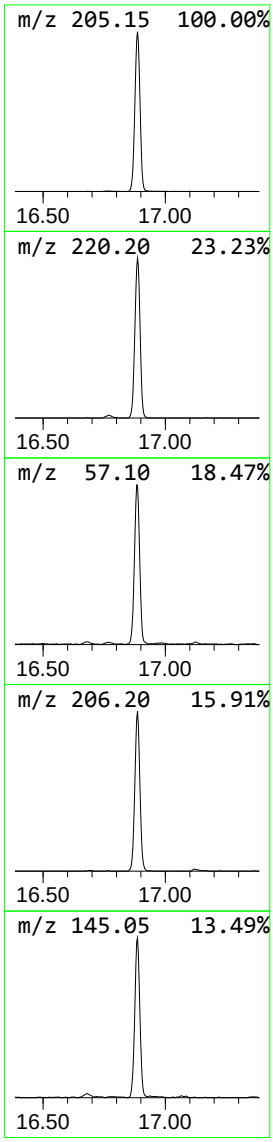
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Butylated Hydroxytoluene Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.886	2.76 ug/L	1556210	1,4-Dichlorobenzene-d4	11.809

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	99
2	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	98
3	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	97
4	Butylated Hydroxytoluene	220	C15H24O	000128-37-0	96
5	Ethyl 4-oxo-2-phenylpentanoate	220	C13H16O3	1000427-11-2	94





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU103122\
Data File : VU051647.D
Acq On : 01 Nov 2022 06:38
Operator : JC/MD
Sample : N5319-07
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 46 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
BHAA0

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR102822WMA.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
Diisopropyl ether	3.990	0.7	ug/L	12678600	1	6.247	91111000	5.0
Benzene, propyl-	10.902	10.3	ug/L	5812520	3	11.809	2821400	5.0
Benzene, 1-chlo...	10.989	7.2	ug/L	4078170	3	11.809	2821400	5.0
Benzene, 1-ethy...	11.288	7.6	ug/L	4286880	3	11.809	2821400	5.0
Benzene, 1,2,3-...	11.893	9.6	ug/L	5393180	3	11.809	2821400	5.0
Benzene, 1-prop...	12.095	13.5	ug/L	7603020	3	11.809	2821400	5.0
Cyclohexanone, ...	12.475	4.6	ug/L	2571900	3	11.809	2821400	5.0
Benzene, 1-ethy...	12.568	3.1	ug/L	1766320	3	11.809	2821400	5.0
m-Xylene, 5-chl...	12.819	2.7	ug/L	1500140	3	11.809	2821400	5.0
unknown-01	12.857	5.9	ug/L	3325970	3	11.809	2821400	5.0
Benzene, 1,2,4,...	12.970	2.4	ug/L	1345780	3	11.809	2821400	5.0
1-Phenyl-1-butene	13.449	3.3	ug/L	1860270	3	11.809	2821400	5.0
Butylated Hydro...	16.886	2.8	ug/L	1556210	3	11.809	2821400	5.0