

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
 Data File : VU047412.D
 Acq On : 09 Mar 2022 18:21
 Operator : SY/MD
 Sample : N1810-22DL 50X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GBDA7DL

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\SFAMUTR030922WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VU047412.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.604	72	82	92	rBV2	76079	88267	6.14%	0.894%
2	1.919	171	180	194	rBV	42409	58246	4.05%	0.590%
3	1.999	195	205	222	rVB	548747	755736	52.61%	7.653%
4	2.208	261	270	285	rVB	106016	155415	10.82%	1.574%
5	2.324	297	306	316	rVB	15497	22455	1.56%	0.227%
6	2.472	344	352	366	rVB	15761	24564	1.71%	0.249%
7	2.575	372	384	408	rBV2	105684	237161	16.51%	2.402%
8	3.051	519	532	538	rVV2	42205	91773	6.39%	0.929%
9	3.089	539	544	553	rVV2	45005	81697	5.69%	0.827%
10	3.144	553	561	580	rVB3	30666	71552	4.98%	0.725%
11	3.366	616	630	648	rVB	46090	101923	7.09%	1.032%
12	3.581	685	697	713	rVB5	8531	18819	1.31%	0.191%
13	3.703	724	735	748	rVB2	6976	14921	1.04%	0.151%
14	3.838	762	777	788	rBV4	6708	15039	1.05%	0.152%
15	3.922	788	803	811	rVV4	9322	22960	1.60%	0.233%
16	3.986	813	823	837	rVB2	14868	34742	2.42%	0.352%
17	4.099	843	858	872	rBV3	10468	27039	1.88%	0.274%
18	4.186	872	885	900	rVV6	7557	22441	1.56%	0.227%
19	4.340	918	933	946	rVB4	14918	33970	2.36%	0.344%
20	4.536	978	994	1010	rBV	68867	165543	11.52%	1.676%
21	4.642	1010	1027	1061	rVV	35831	126276	8.79%	1.279%
22	5.070	1142	1160	1175	rBV	71344	156026	10.86%	1.580%
23	5.182	1184	1195	1211	rVB3	16528	34475	2.40%	0.349%
24	5.395	1241	1261	1274	rBV	36660	85215	5.93%	0.863%
25	5.465	1274	1283	1297	rVB6	6588	15663	1.09%	0.159%
26	5.729	1345	1365	1375	rBV2	152513	387937	27.00%	3.929%
27	6.250	1512	1527	1552	rBV	198479	402390	28.01%	4.075%
28	6.694	1649	1665	1678	rBV	89052	175443	12.21%	1.777%
29	6.764	1681	1687	1699	rVB5	8626	15985	1.11%	0.162%
30	7.395	1871	1883	1895	rBV4	8055	14645	1.02%	0.148%
31	7.568	1925	1937	1956	rBV	50005	87908	6.12%	0.890%
32	7.899	2026	2040	2052	rBV	204253	349677	24.34%	3.541%
33	7.970	2052	2062	2088	rVB	184758	322962	22.48%	3.271%
34	8.182	2117	2128	2143	rBV	30143	52054	3.62%	0.527%
35	8.475	2207	2219	2235	rVB2	8034	14789	1.03%	0.150%
36	8.636	2256	2269	2310	rBV	186591	393703	27.41%	3.987%

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37	9.417	2498	2512	2540	rBV	279099	462764	32.21%	4.686%
38	9.571	2547	2560	2584	rBV	315589	509429	35.46%	5.159%
39	9.694	2584	2598	2629	rVB	872045	1436598	100.00%	14.549%
40	10.102	2709	2725	2750	rBV	203880	327309	22.78%	3.315%
41	10.484	2834	2844	2857	rVB	17914	26826	1.87%	0.272%
42	10.758	2915	2929	2947	rBV	71785	116330	8.10%	1.178%
43	10.906	2956	2975	2989	rBV2	61460	105329	7.33%	1.067%
44	10.999	2992	3004	3022	rVV	137996	304820	21.22%	3.087%
45	11.089	3022	3032	3046	rVB	64334	98552	6.86%	0.998%
46	11.292	3081	3095	3110	rBV	70205	110315	7.68%	1.117%
47	11.468	3137	3150	3165	rBV	279129	423750	29.50%	4.291%
48	11.812	3244	3257	3272	rBV	327305	515515	35.88%	5.221%
49	11.896	3272	3283	3297	rVB	63678	98933	6.89%	1.002%
50	12.095	3330	3345	3363	rBV2	74112	127504	8.88%	1.291%
51	12.192	3363	3375	3392	rVB	211014	345410	24.04%	3.498%
52	12.571	3482	3493	3500	rBV	11093	15675	1.09%	0.159%
53	12.681	3517	3527	3542	rBV3	15758	25934	1.81%	0.263%
54	12.973	3605	3618	3626	rBV	12562	19242	1.34%	0.195%
55	13.028	3626	3635	3651	rVB	13436	20546	1.43%	0.208%
56	13.295	3708	3718	3734	rVB	13448	19791	1.38%	0.200%
57	13.455	3756	3768	3778	rVB3	24326	39086	2.72%	0.396%
58	14.089	3953	3965	3976	rBV	28190	45378	3.16%	0.460%

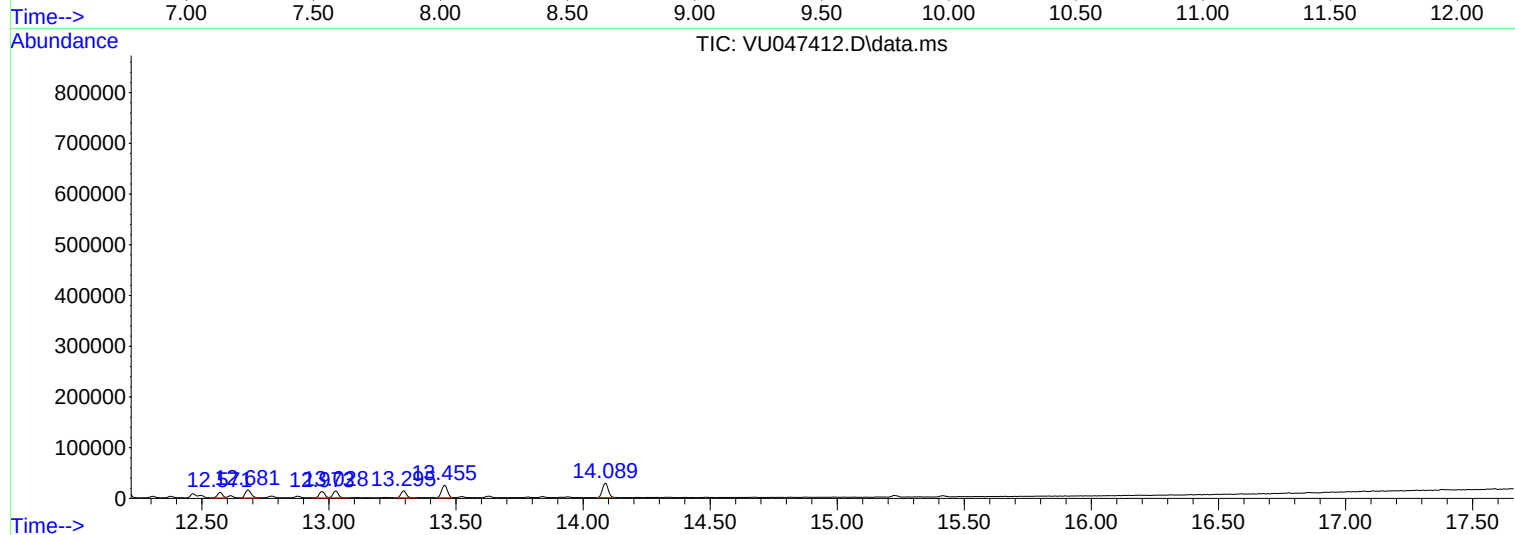
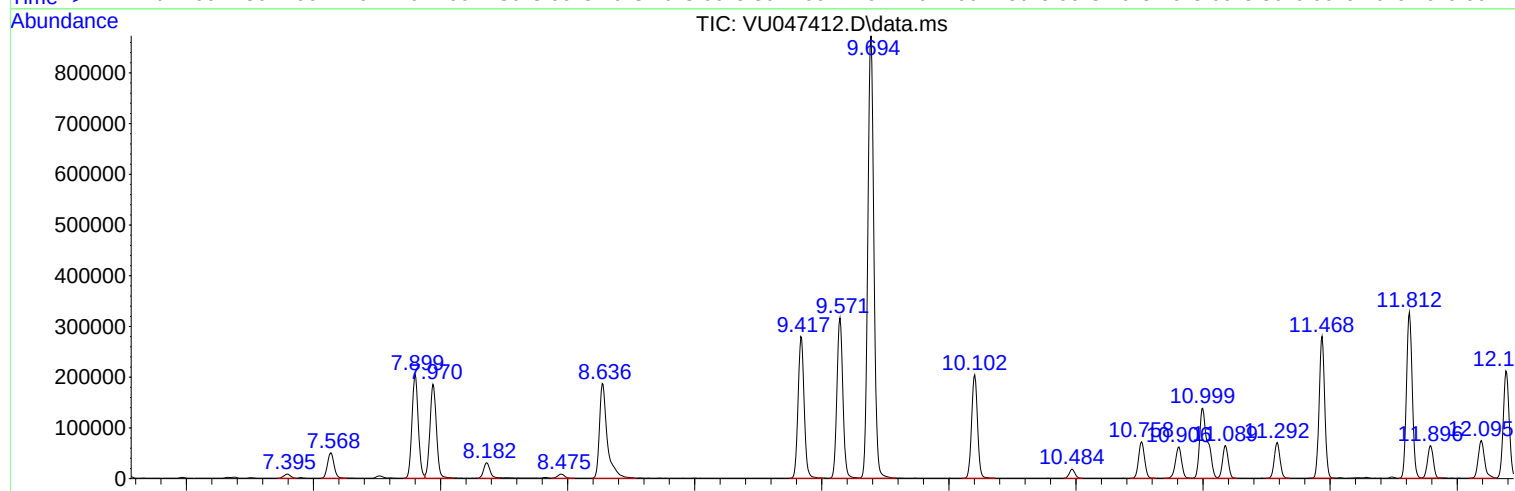
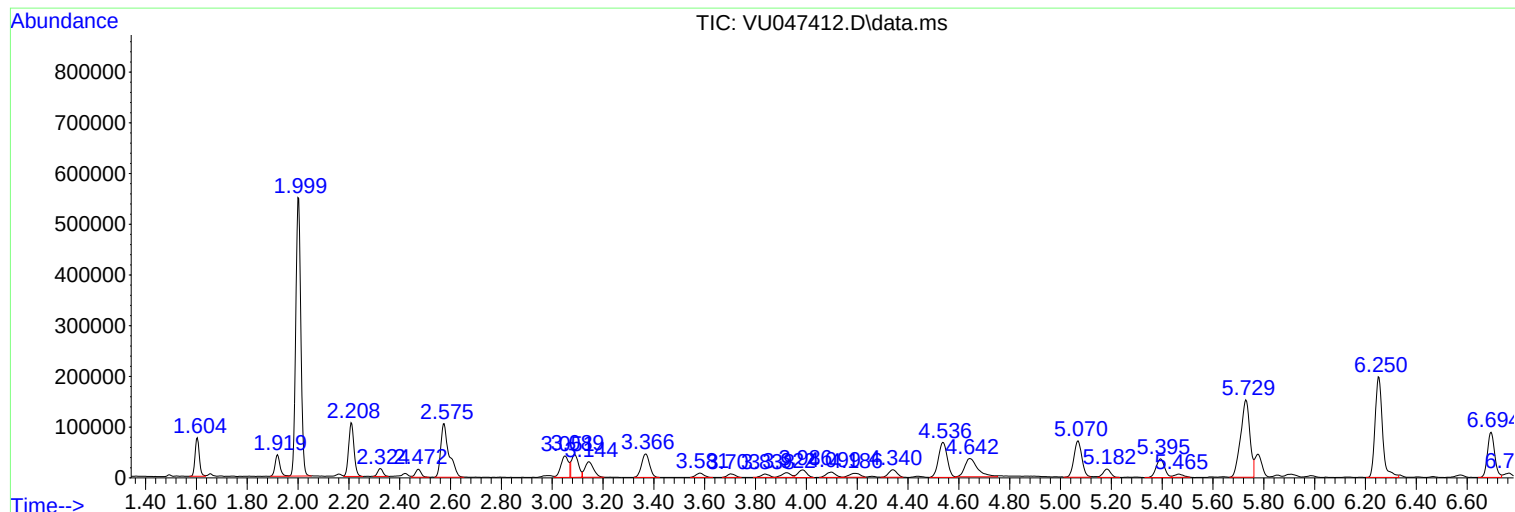
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Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR030922WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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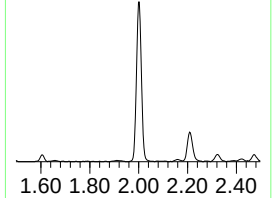
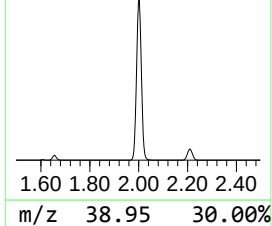
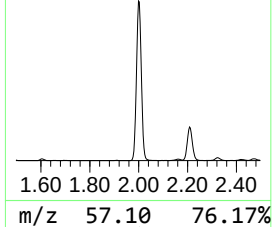
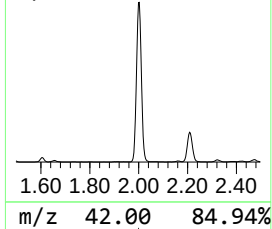
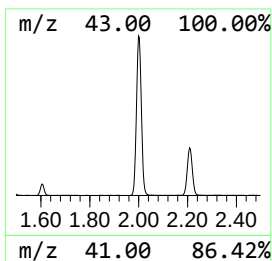
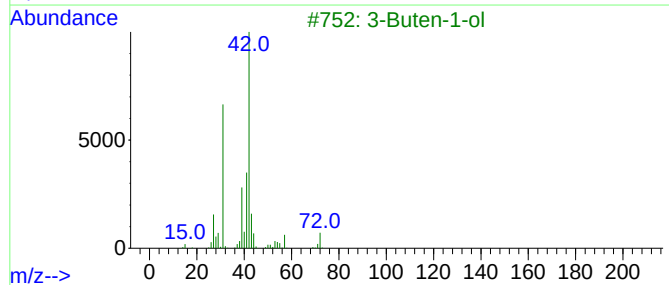
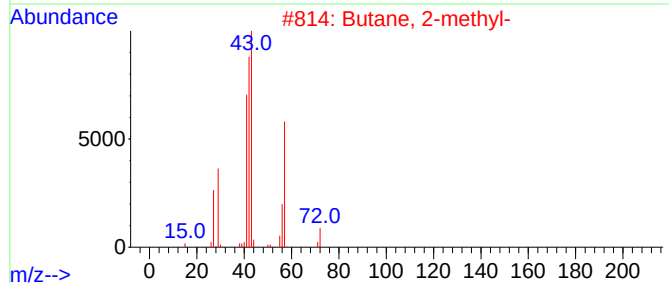
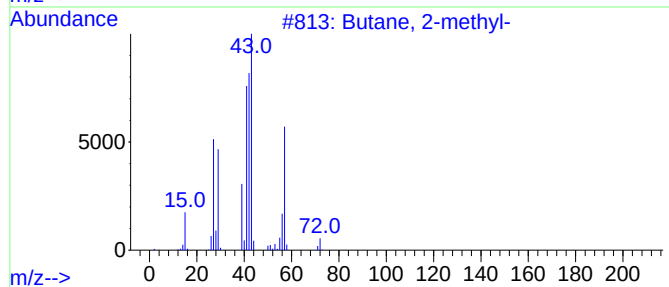
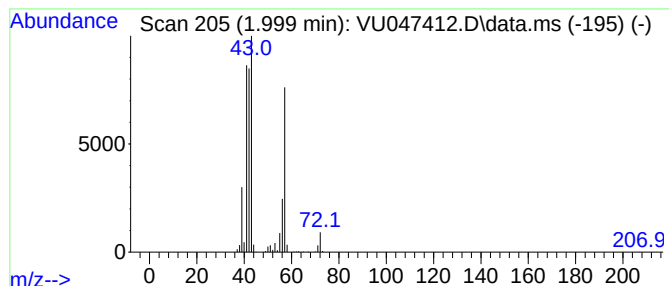
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Quant Title : TRACE VOA SFAM1.0

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.999	9.39 ug/L	755736	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-			72	C5H12	000078-78-4	90
2	Butane, 2-methyl-			72	C5H12	000078-78-4	87
3	3-Buten-1-ol			72	C4H8O	000627-27-0	43
4	Pentane			72	C5H12	000109-66-0	42
5	Pentane			72	C5H12	000109-66-0	28



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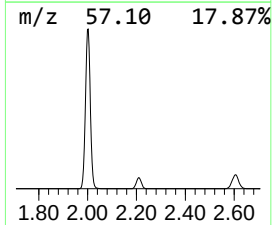
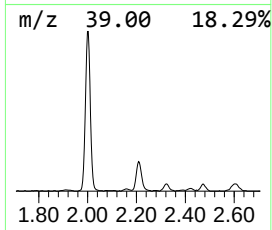
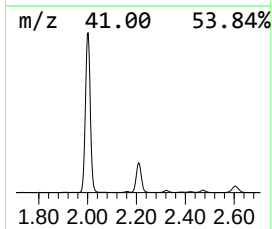
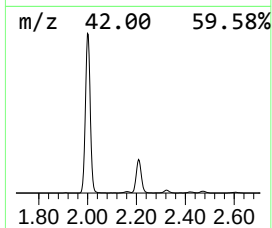
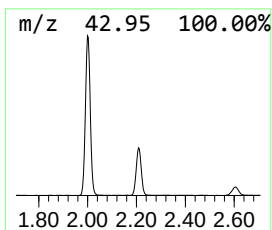
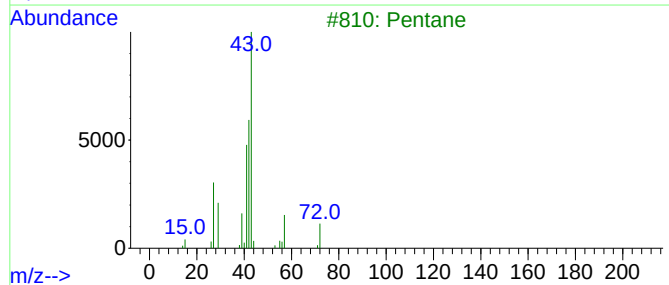
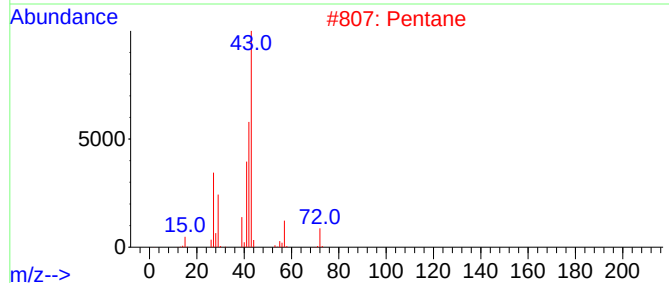
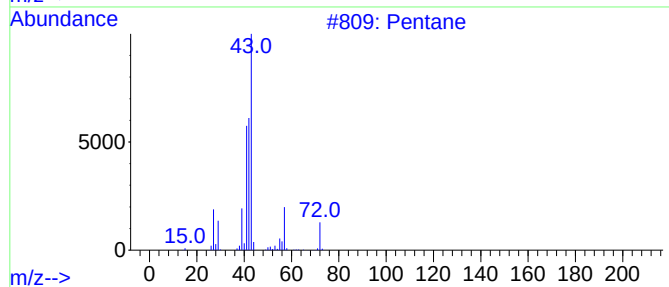
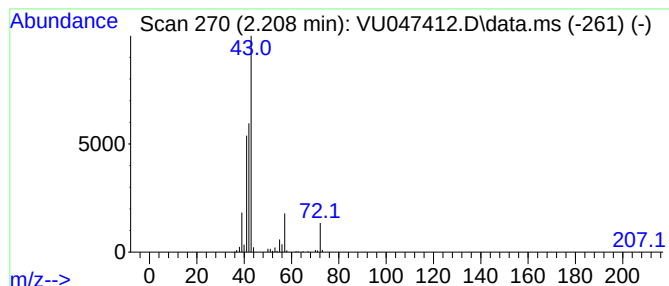
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR030922WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.208	1.93 ug/L	155415	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane			72	C5H12	000109-66-0	91
2	Pentane			72	C5H12	000109-66-0	90
3	Pentane			72	C5H12	000109-66-0	90
4	Pentane			72	C5H12	000109-66-0	90
5	Pentane			72	C5H12	000109-66-0	90



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Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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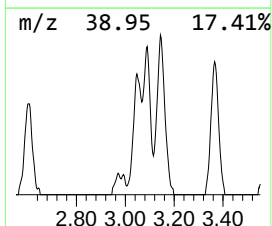
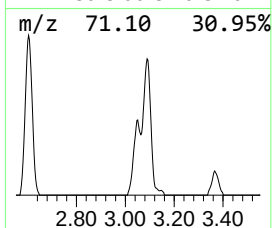
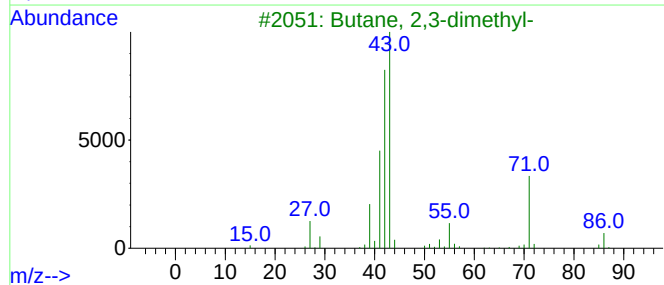
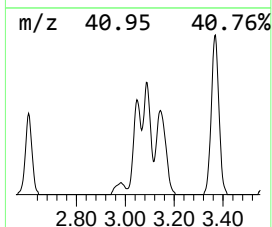
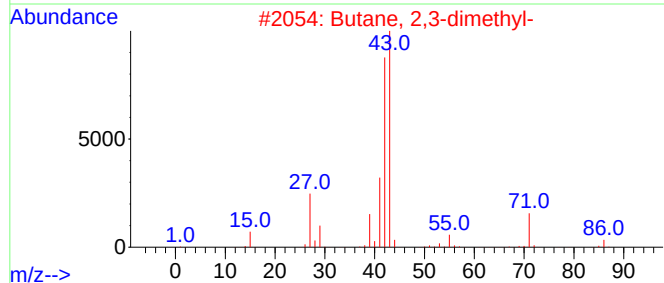
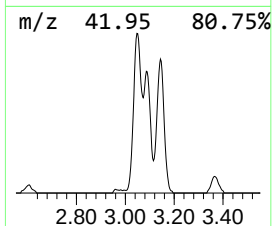
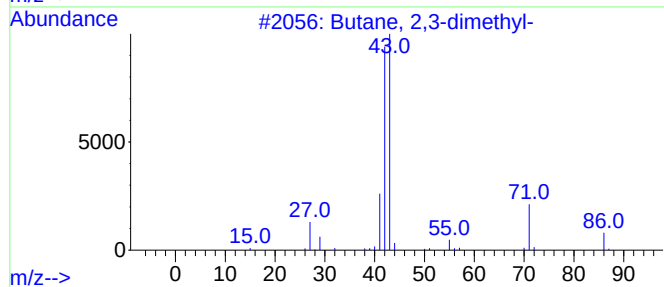
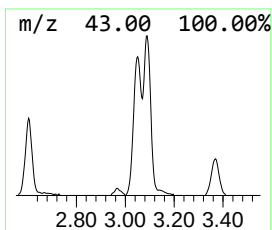
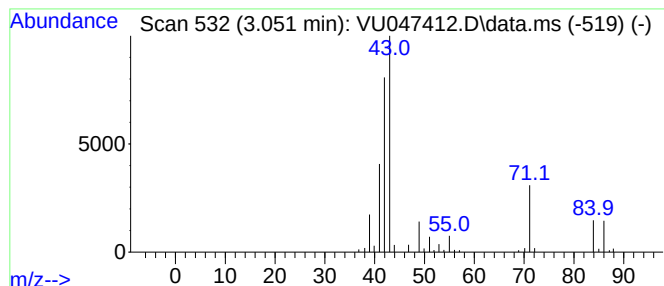
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.051	1.14 ug/L	91773	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64
2			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	64
3			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	59
4			Butane, 2,3-dimethyl-	86	C6H14	000079-29-8	53
5			Pentane, 2-methyl-	86	C6H14	000107-83-5	38



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Instrument :
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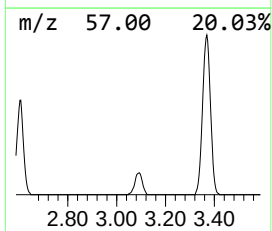
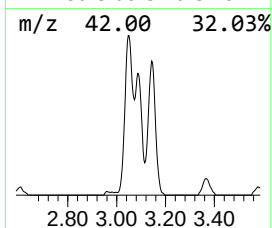
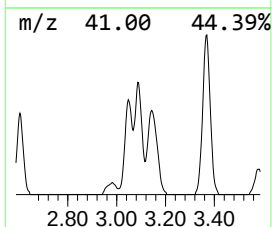
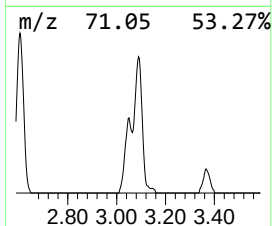
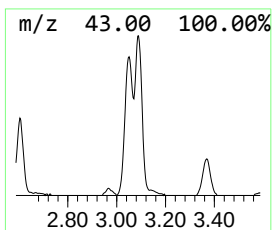
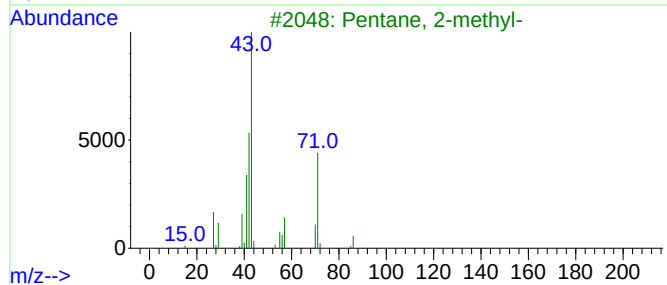
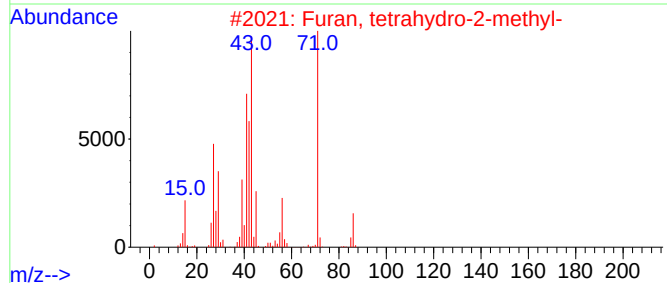
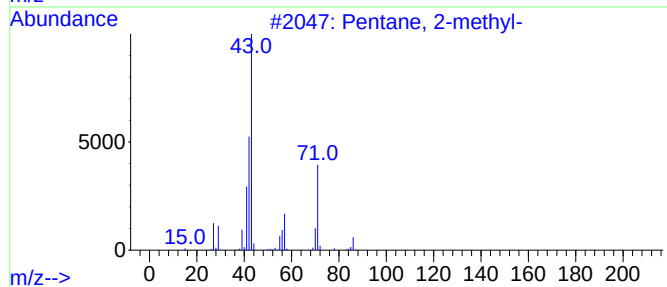
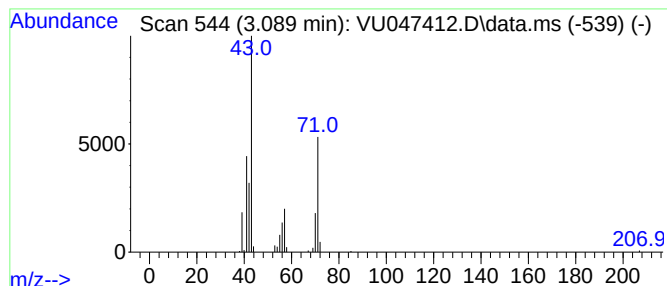
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.089	1.02 ug/L	81697	1,4-Difluorobenzene	6.250

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentane, 2-methyl-	86	C6H14	000107-83-5	53
2			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	50
3			Pentane, 2-methyl-	86	C6H14	000107-83-5	45
4			Pentane, 3,3-dimethyl-	100	C7H16	000562-49-2	42
5			Furan, tetrahydro-2-methyl-	86	C5H10O	000096-47-9	38



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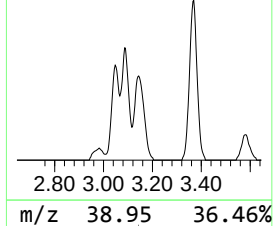
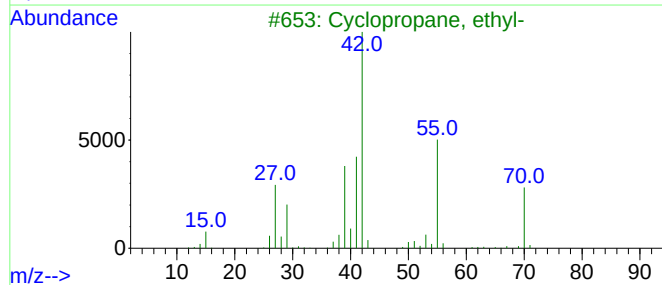
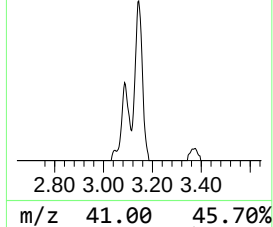
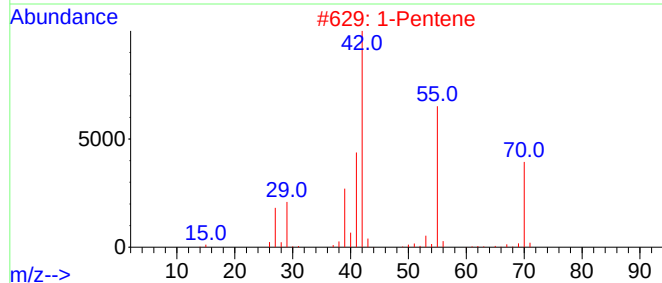
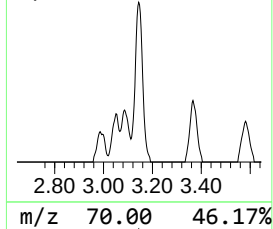
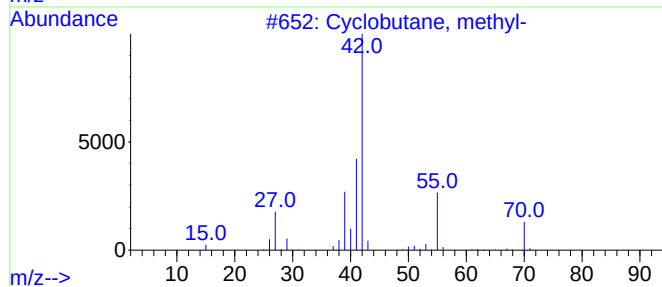
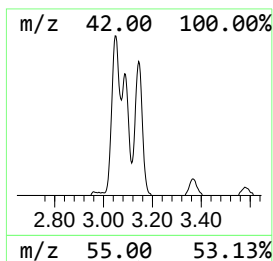
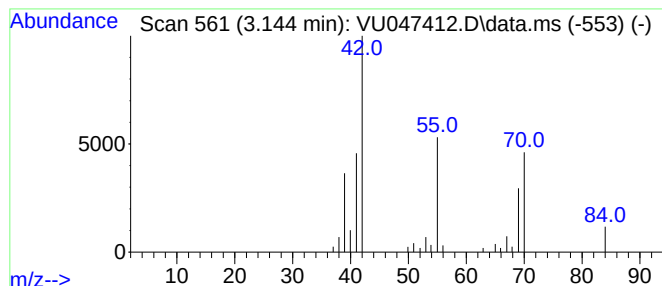
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 5 (DEL) Alkane: Cyclic3.144 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.144	0.89 ug/L	71552	1,4-Difluorobenzene	6.250

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclobutane, methyl-	70	C5H10	000598-61-8	72
2	1-Pentene	70	C5H10	000109-67-1	72
3	Cyclopropane, ethyl-	70	C5H10	001191-96-4	72
4	Cyclopropane, ethyl-	70	C5H10	001191-96-4	64
5	Cyclopentane	70	C5H10	000287-92-3	50



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

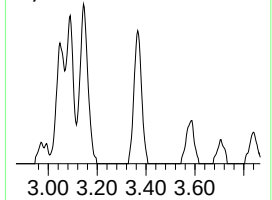
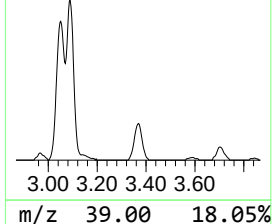
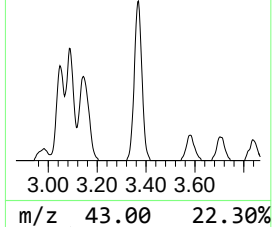
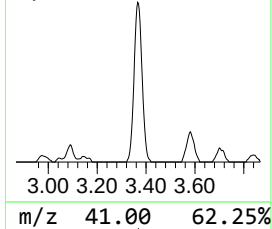
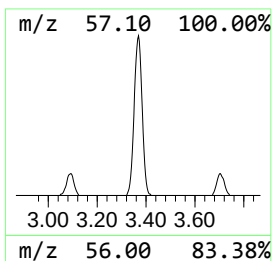
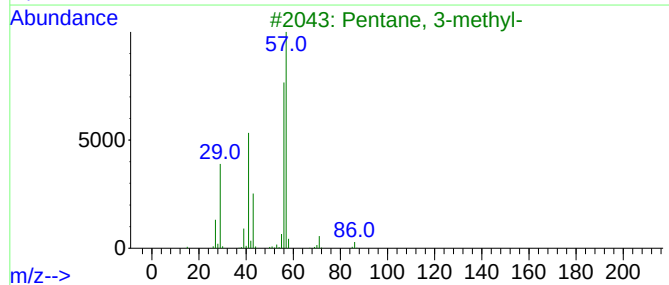
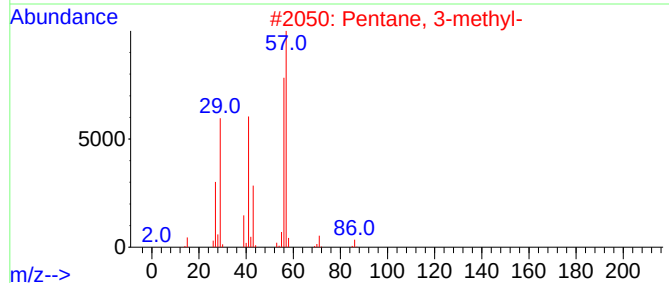
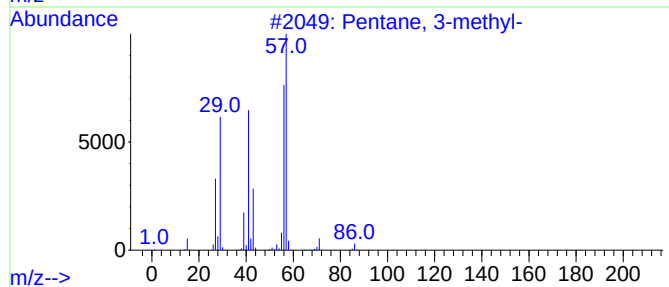
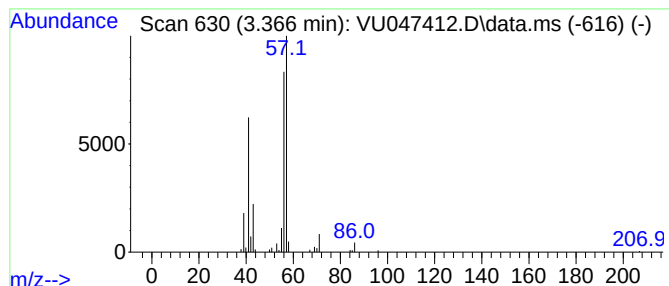
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.366	1.27 ug/L	101923	1,4-Difluorobenzene	6.250

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2		Pentane, 3-methyl-	86	C6H14	000096-14-0	90
3		Pentane, 3-methyl-	86	C6H14	000096-14-0	86
4		Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	78
5		Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

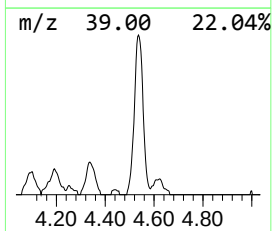
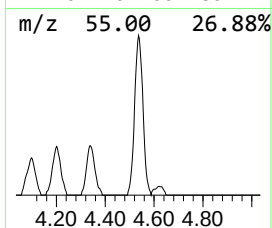
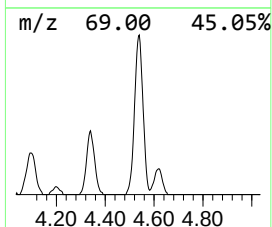
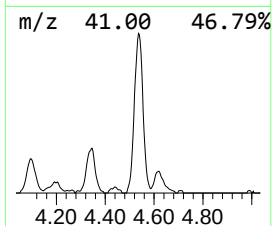
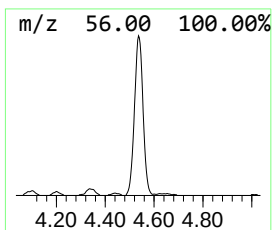
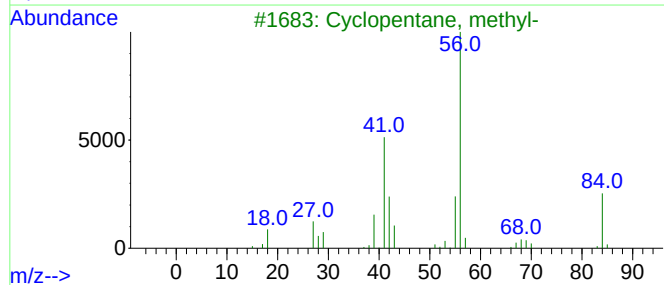
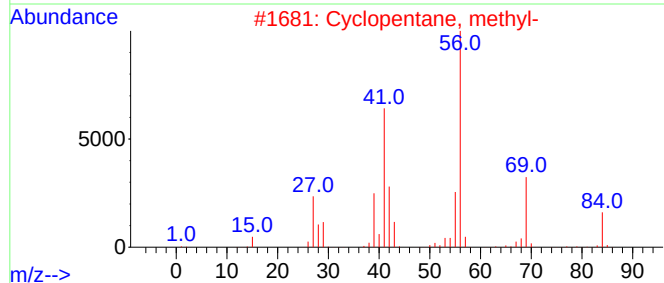
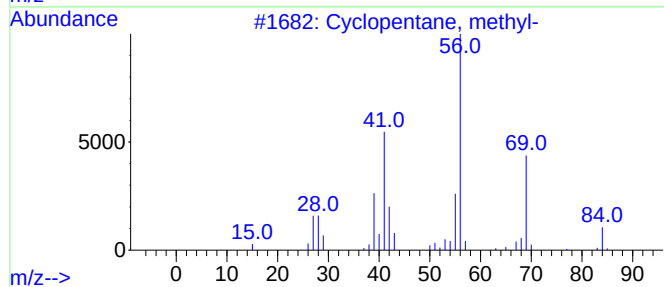
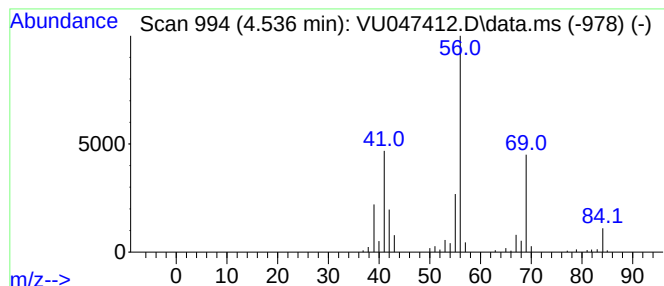
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 7 (DEL) Alkane: Cyclic4.536 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.536	2.06 ug/L	165543	1,4-Difluorobenzene	6.250

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclopentane, methyl-	84	C6H12	000096-37-7	91
2		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
3		Cyclopentane, methyl-	84	C6H12	000096-37-7	86
4		1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	80
5		Cyclopentane, methyl-	84	C6H12	000096-37-7	78



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

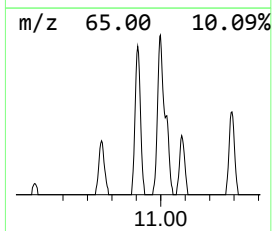
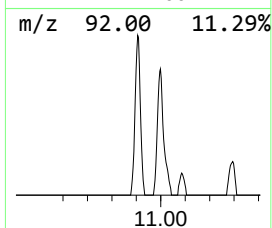
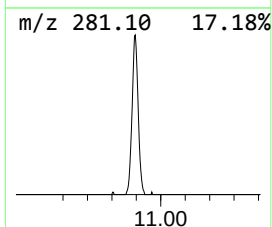
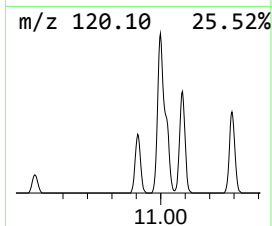
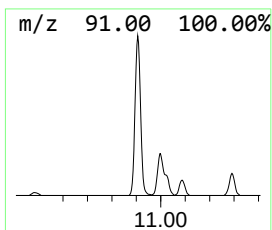
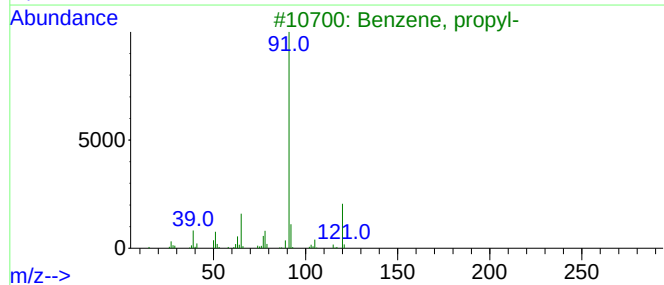
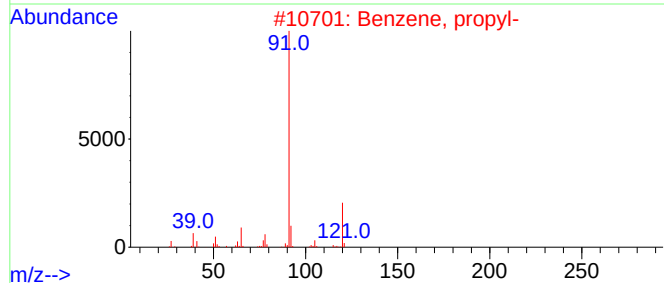
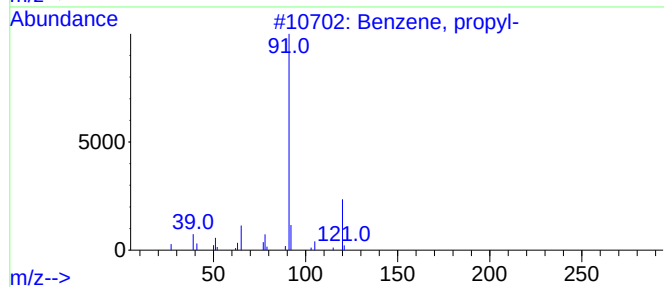
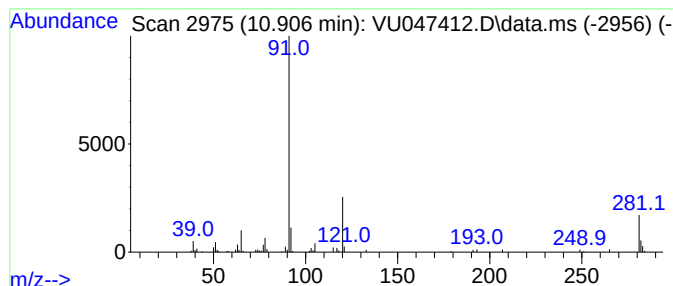
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 9 Benzene, propyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.906	1.02 ug/L	105329	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	92
2			Benzene, propyl-	120	C9H12	000103-65-1	70
3			Benzene, propyl-	120	C9H12	000103-65-1	70
4			Benzene, propyl-	120	C9H12	000103-65-1	70
5			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

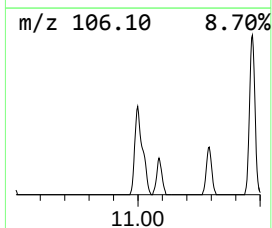
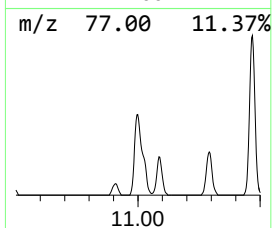
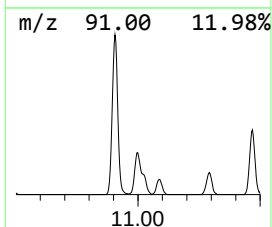
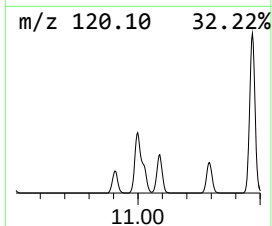
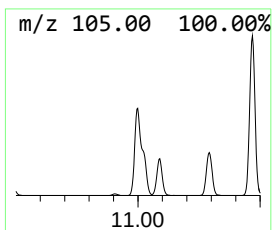
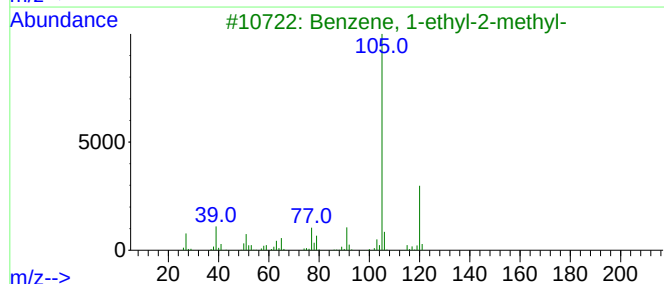
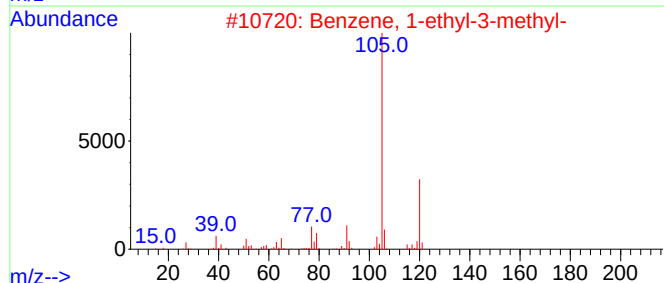
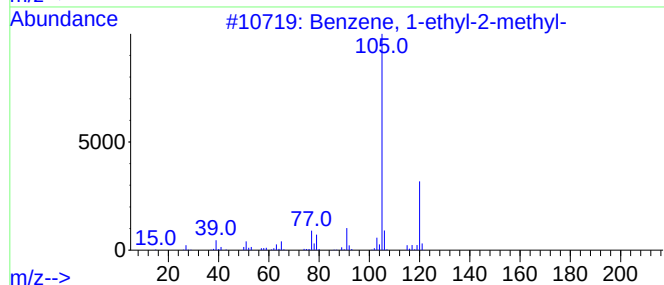
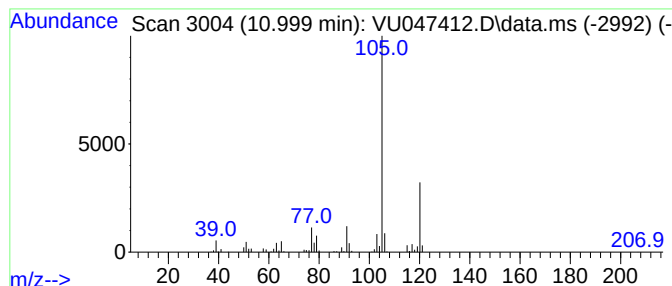
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 10 Benzene, 1-ethyl-2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.999	2.96 ug/L	304820	1,4-Dichlorobenzene-d4	11.812

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			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

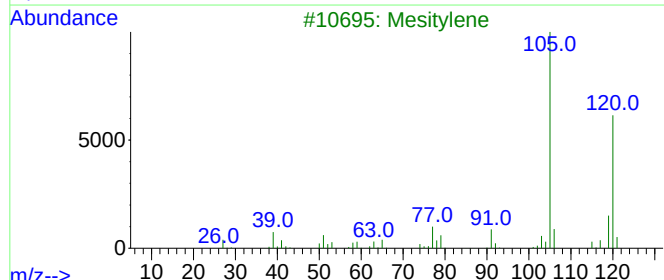
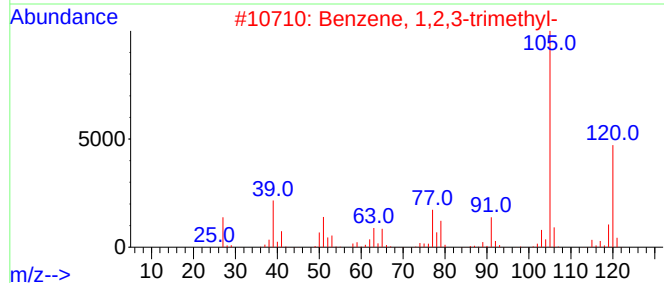
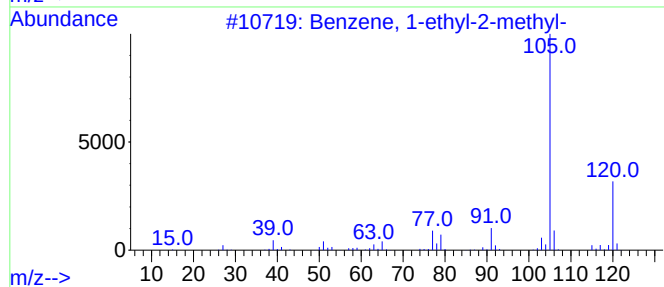
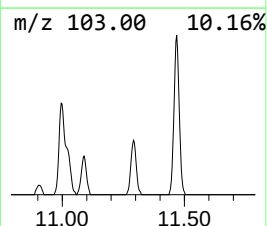
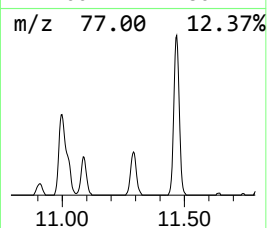
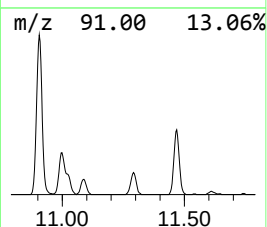
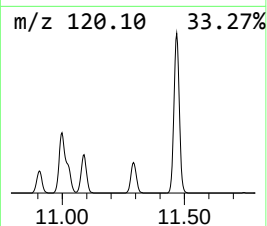
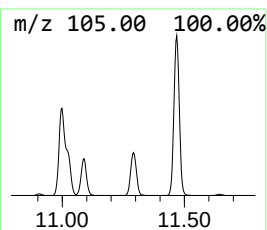
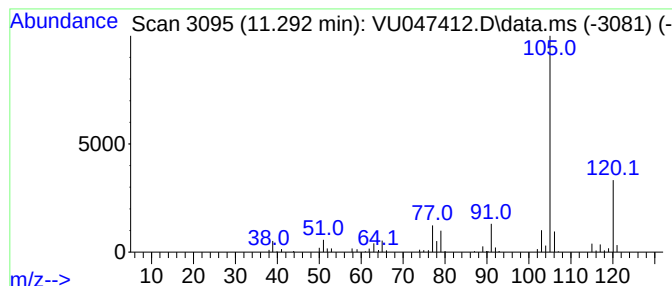
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 11 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.291	1.07 ug/L	110315	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3			Mesitylene	120	C9H12	000108-67-8	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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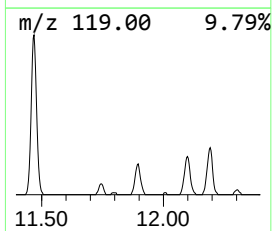
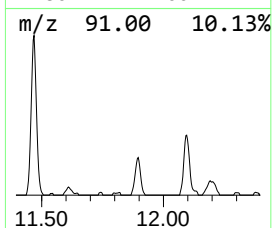
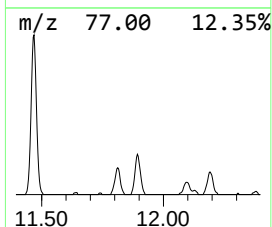
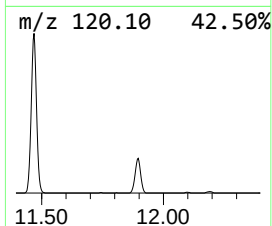
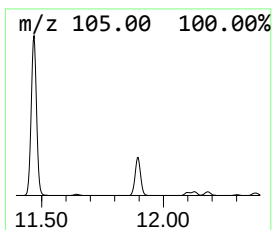
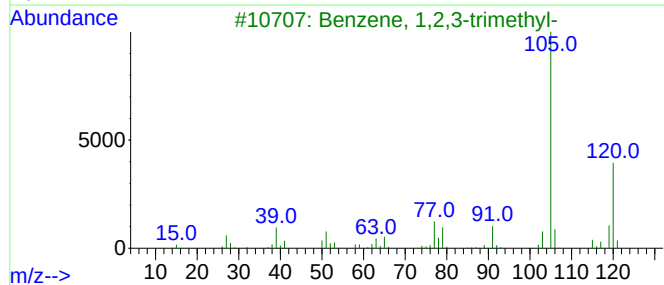
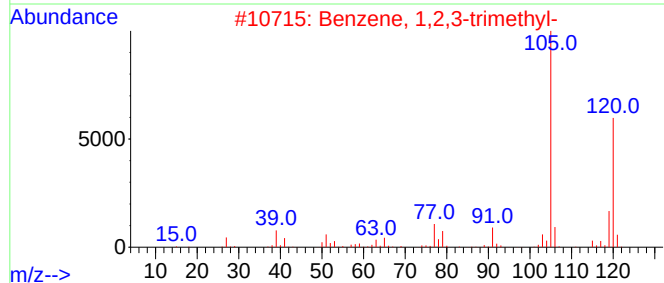
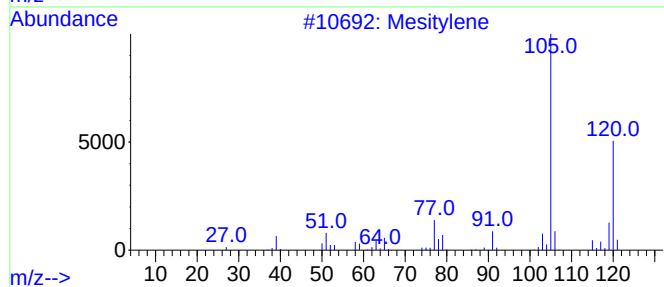
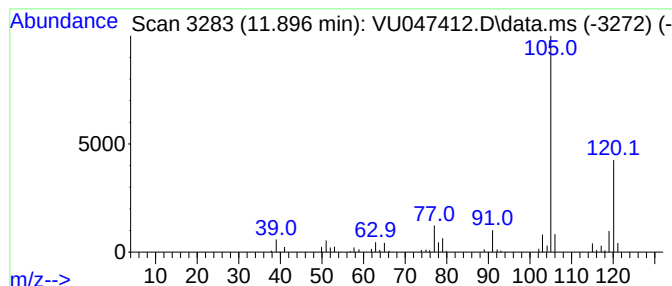
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 12 unknown-01 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.896	0.96 ug/L	98933	1,4-Dichlorobenzene-d4	11.812

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	94
4			Mesitylene	120	C9H12	000108-67-8	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
Data File : VU047412.D
Acq On : 09 Mar 2022 18:21
Operator : SY/MD
Sample : N1810-22DL 50X
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
GBDA7DL

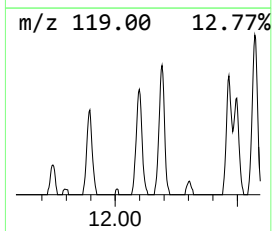
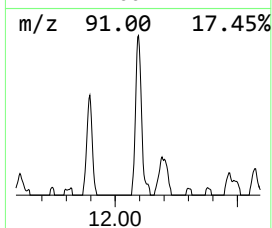
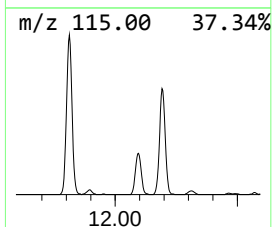
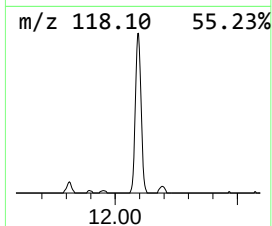
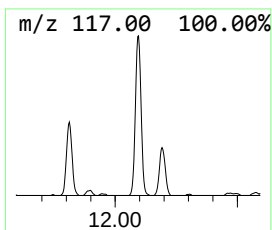
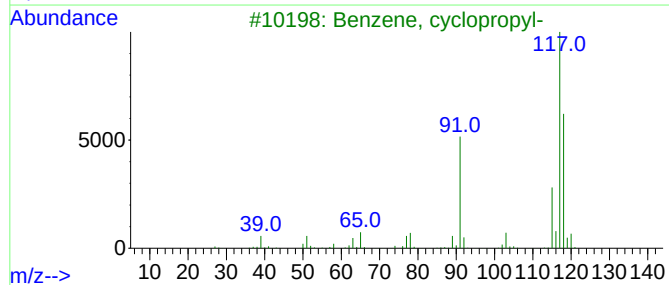
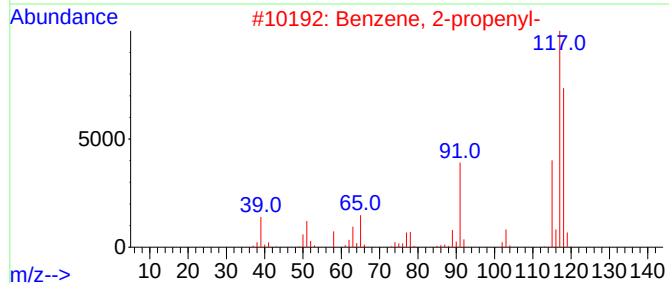
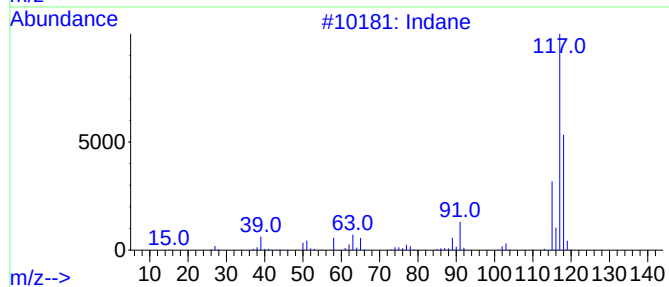
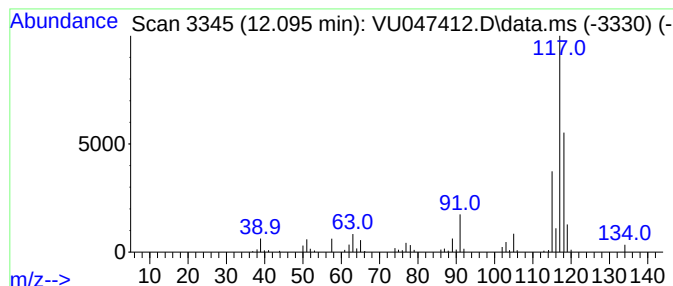
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 13 Indane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.095	1.24 ug/L	127504	1,4-Dichlorobenzene-d4	11.812

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	74
3			Benzene, cyclopropyl-	118	C9H10	000873-49-4	74
4			Deltacyclene	118	C9H10	007785-10-6	64
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU030922\
 Data File : VU047412.D
 Acq On : 09 Mar 2022 18:21
 Operator : SY/MD
 Sample : N1810-22DL 50X
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 GBDA7DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR030922WMA.M
 Quant Title : TRACE VOA SFAM1.0

TIC Library : C:\Database\NIST0.L
 TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: S...	1.999	9.4	ug/L	755736	1	6.250	402390	5.0
(DEL) Alkane: S...	2.208	1.9	ug/L	155415	1	6.250	402390	5.0
(DEL) Alkane: S...	3.051	1.1	ug/L	91773	1	6.250	402390	5.0
(DEL) Alkane: S...	3.089	1.0	ug/L	81697	1	6.250	402390	5.0
(DEL) Alkane: C...	3.144	0.9	ug/L	71552	1	6.250	402390	5.0
(DEL) Alkane: S...	3.366	1.3	ug/L	101923	1	6.250	402390	5.0
(DEL) Alkane: C...	4.536	2.1	ug/L	165543	1	6.250	402390	5.0
Benzene, propyl-	10.906	1.0	ug/L	105329	3	11.812	515515	5.0
Benzene, 1-ethy...	10.999	3.0	ug/L	304820	3	11.812	515515	5.0
Benzene, 1,2,3-...	11.291	1.1	ug/L	110315	3	11.812	515515	5.0
unknown-01	11.896	1.0	ug/L	98933	3	11.812	515515	5.0
Indane	12.095	1.2	ug/L	127504	3	11.812	515515	5.0