LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MDL-WATER-QT3-2021-06

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\SFAMUTR051521WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VU043793.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	73	82	95	rBV2	178025	213514	6.25%	0.898%
2	1.919	171	180	193	rVB	78988	103541	3.03%	0.435%
3	2.572	369	383	402	rBV2	182781	316582	9.26%	1.331%
4	3.054	516	533	535	rBV4	184397	310839	9.09%	1.307%
5	3.089	537	544	569	rVB	465259	924751	27.06%	3.889%
6	3.369	611	631	657	rBV	1307544	2895441	84.71%	12.177%
7	3.707	717	736	761	rBV	836114	1847491	54.05%	7.770%
8	3.880	765	790	813	rBV	66026	207582	6.07%	0.873%
9	3.986	814	823	840	rVB3	23477	54858	1.60%	0.231%
10	4.099	844	858	872	rBV4	21397	49860	1.46%	0.210%
11	4.308	903	923	948	rBV6	45621	181095	5.30%	0.762%
12	4.443	951	965	977	rBV2	40250	94206	2.76%	0.396%
13	4.540	977	995	1013	rBV	1425595	3417947	100.00%	14.374%
14	4.675	1014	1037	1065	rVB2	220434	712005	20.83%	2.994%
15	5.073	1145	1161	1178	rBV	141970	316982	9.27%	1.333%
16	5.324	1220	1239	1251	rBV	187566	422628	12.36%	1.777%
17	5.398	1251	1262	1279	rVB2	150564	335204	9.81%	1.410%
18	5.736	1345	1367	1395	rBV2	304775	785005	22.97%	3.301%
19	6.256	1513	1529	1548	rBV	284825	534146	15.63%	2.246%
20	6.700	1653	1667	1684	rBV	178696	350400	10.25%	1.474%
21	7.572	1918	1938	1954	rBV	88024	155315	4.54%	0.653%
22	7.906	2026	2042	2053	rBV	363075	616618	18.04%	2.593%
23	7.977	2053	2064	2082	rVB	544754	931870	27.26%	3.919%
24	8.186	2117	2129	2143	rBV	55002	92648	2.71%	0.390%
25	8.642	2258	2271	2304	rBV	389221	790347	23.12%	3.324%
26	9.424	2500	2514	2521	rBV	410051	748415	21.90%	3.147%
27	9.578	2550	2562	2581	rVB	123549	202253	5.92%	0.851%
28	9.700	2585	2600	2621	rBV	203285	337321	9.87%	1.419%
29	10.108	2714	2727	2744	rBV	102556	164927	4.83%	0.694%
30	10.491	2835	2846	2857	rBV	44679	67531	1.98%	0.284%
31	10.764	2916	2931	2950	rBV	147308	230603	6.75%	0.970%
32	10.912	2959	2977	2992	rBV	126781	192987	5.65%	0.812%
33	11.005	2992	3006	3010	rBV	250229	396760	11.61%	1.669%
34	11.096	3024	3034	3049	rVB	151382	232612	6.81%	0.978%
35	11.240	3065	3079	3088	rBV	446655	697272	20.40%	2.932%
36	11.298	3088	3097	3114	rVB	248279	395468	11.57%	1.663%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MDL-WATER-QT3-2021-06

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\SFAMUTR051521WMA.M
Title : TRACE VOA SFAM1.0

37	11.475	3139	3152	3169	rBV	706561	1079892	31.59%	4.541%
38	11.565	3169	3180	3191	rBV	45275	72937	2.13%	0.307%
39	11.751	3228	3238	3247	rBV	56483	86969	2.54%	0.366%
40	11.819	3247	3259	3275	rVV2	472354	875505	25.61%	3.682%
41	11.899	3275	3284	3297	rVB	154464	236741	6.93%	0.996%
42	12.102	3334	3347	3364	rBV	122398	206364	6.04%	0.868%
43	12.198	3364	3377	3396	rVB	398659	693087	20.28%	2.915%
44	12.578	3484	3495	3504	rBV	26387	39738	1.16%	0.167%
45	13.031	3628	3636	3648	rVB	27648	40863	1.20%	0.172%
46	13.459	3757	3769	3784	rVB3	29038	49696	1.45%	0.209%
47	14.092	3955	3966	3979	rBV	43333	69641	2.04%	0.293%

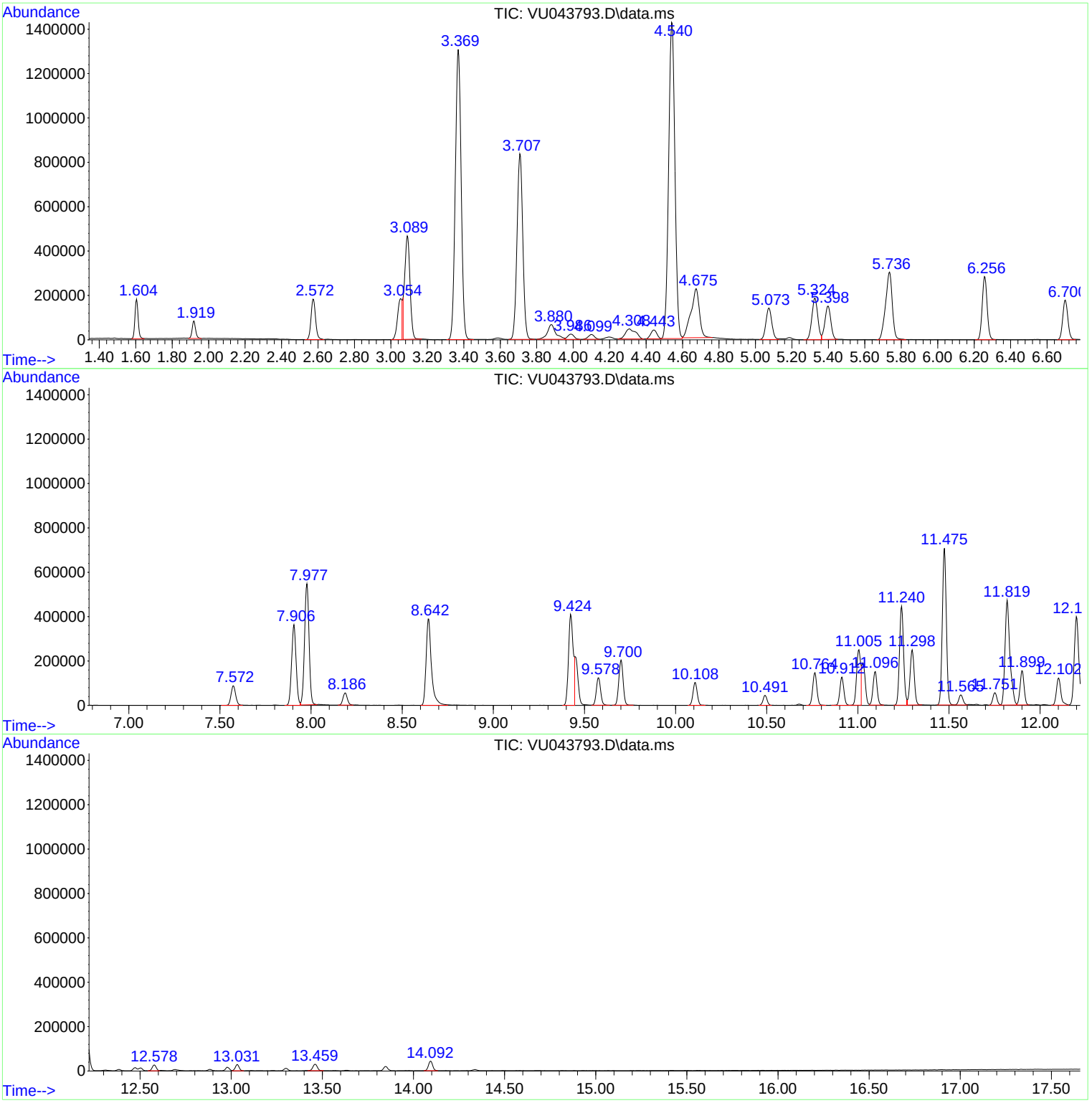
Sum of corrected areas: 23778457

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

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



Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

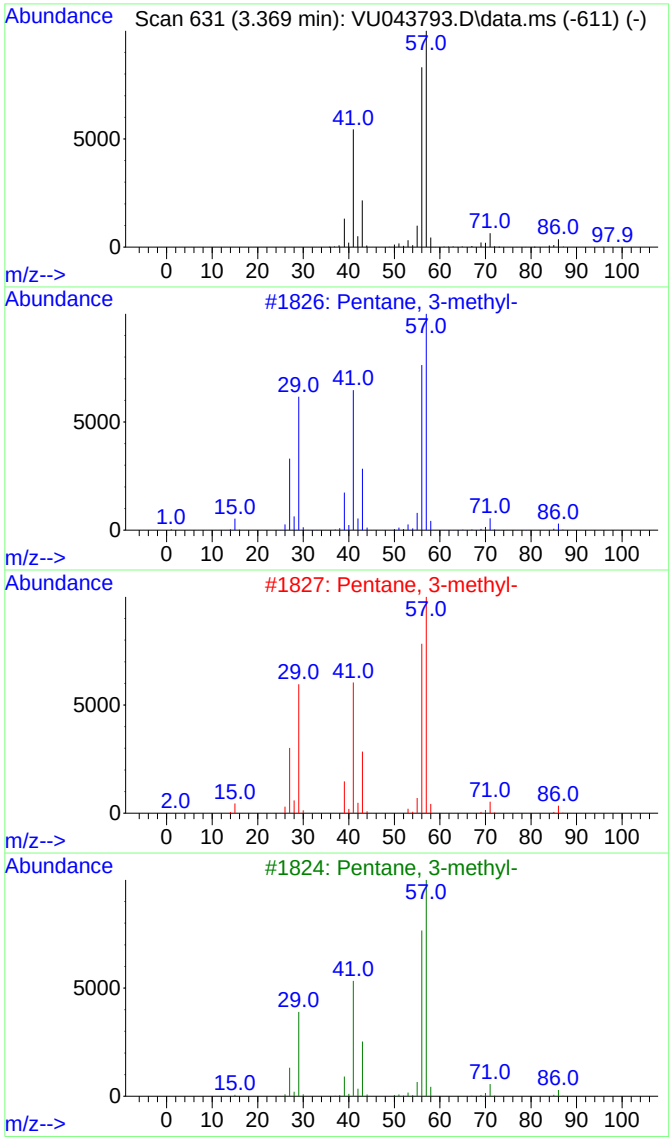
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

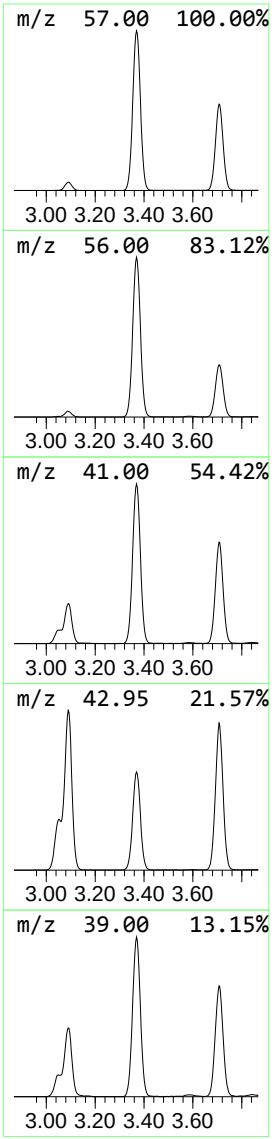
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.369	27.10 ug/L	2895440	1,4-Difluorobenzene	6.256

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	91
3	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
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\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

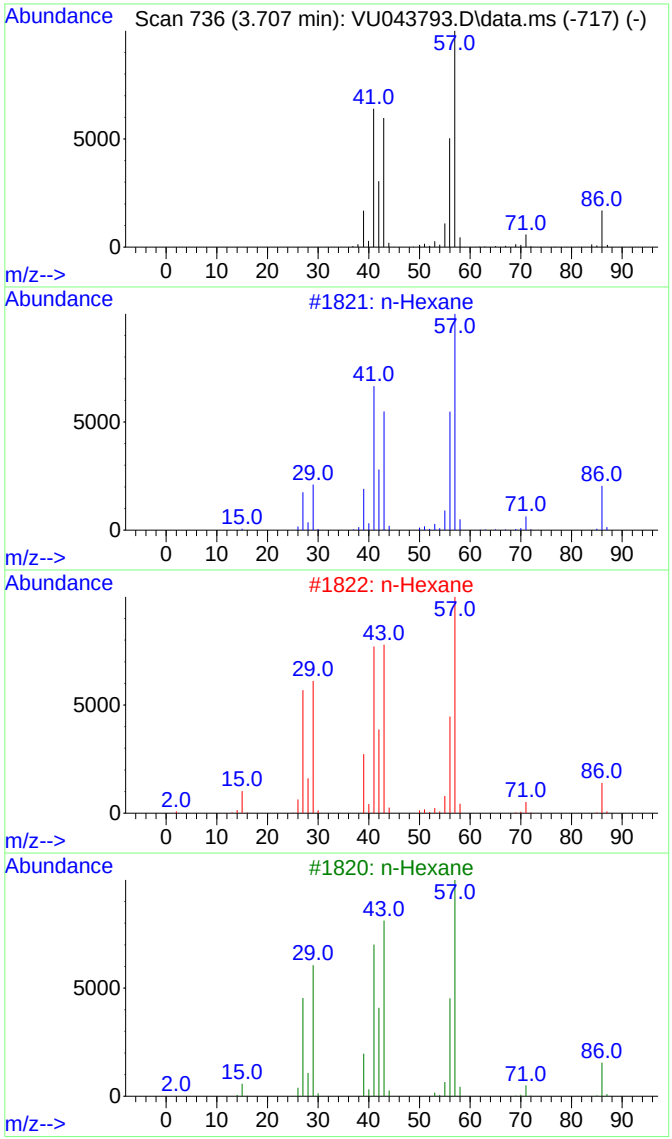
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

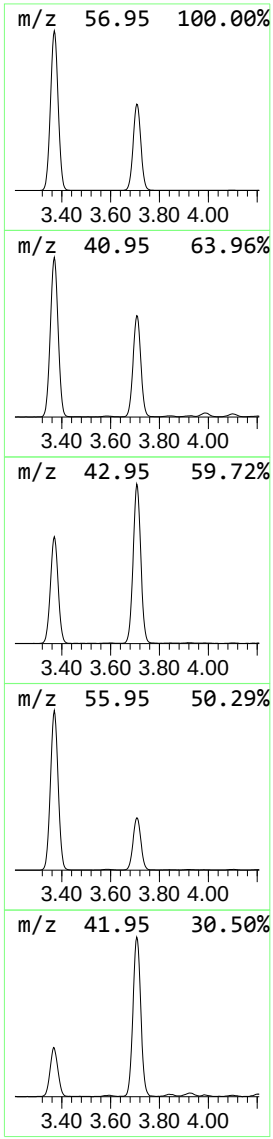
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.707	17.29 ug/L	1847490	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexane	86	C6H14	000110-54-3	94
2	n-Hexane	86	C6H14	000110-54-3	83
3	n-Hexane	86	C6H14	000110-54-3	72
4	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	53
5	Furan, tetrahydro-3-methyl-	86	C5H10O	013423-15-9	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

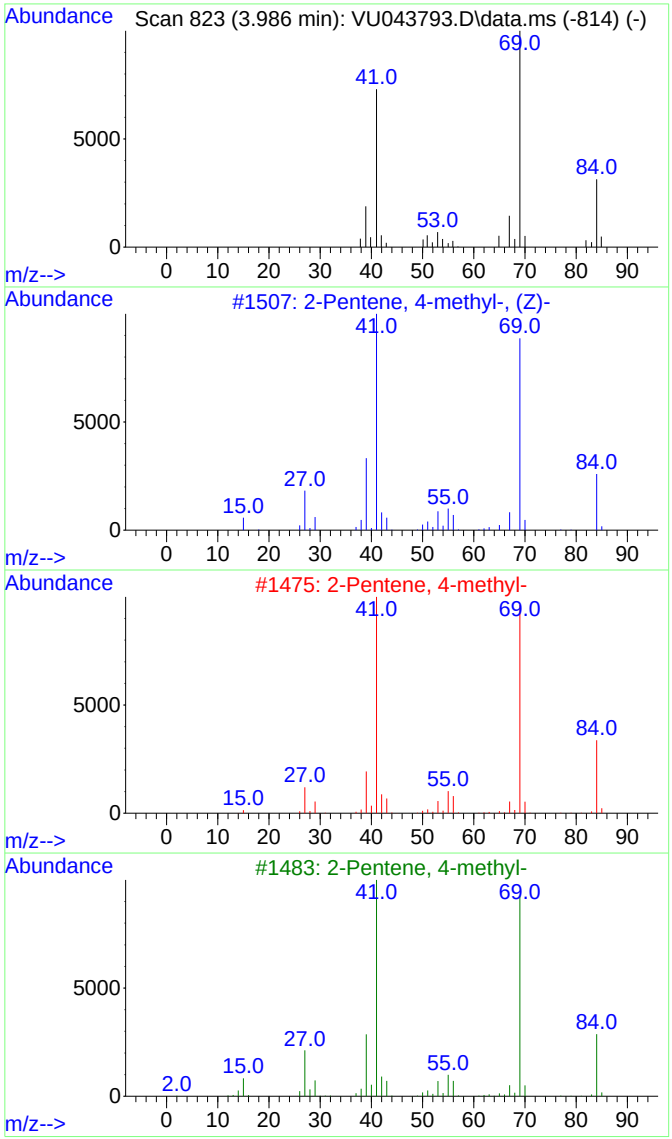
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

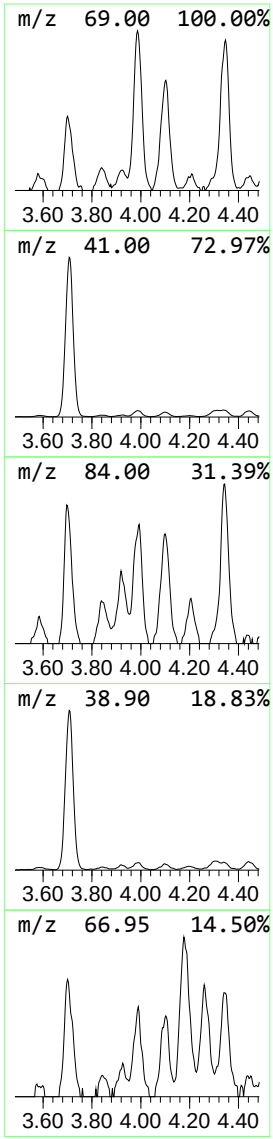
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.986	0.51 ug/L	54858	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentene, 4-methyl-, (Z)-	84	C6H12	000691-38-3	86
2	2-Pentene, 4-methyl-	84	C6H12	004461-48-7	80
3	2-Pentene, 4-methyl-	84	C6H12	004461-48-7	80
4	2-Pentene, 4-methyl-, (E)-	84	C6H12	000674-76-0	78
5	2-Butene, 2,3-dimethyl-	84	C6H12	000563-79-1	78





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

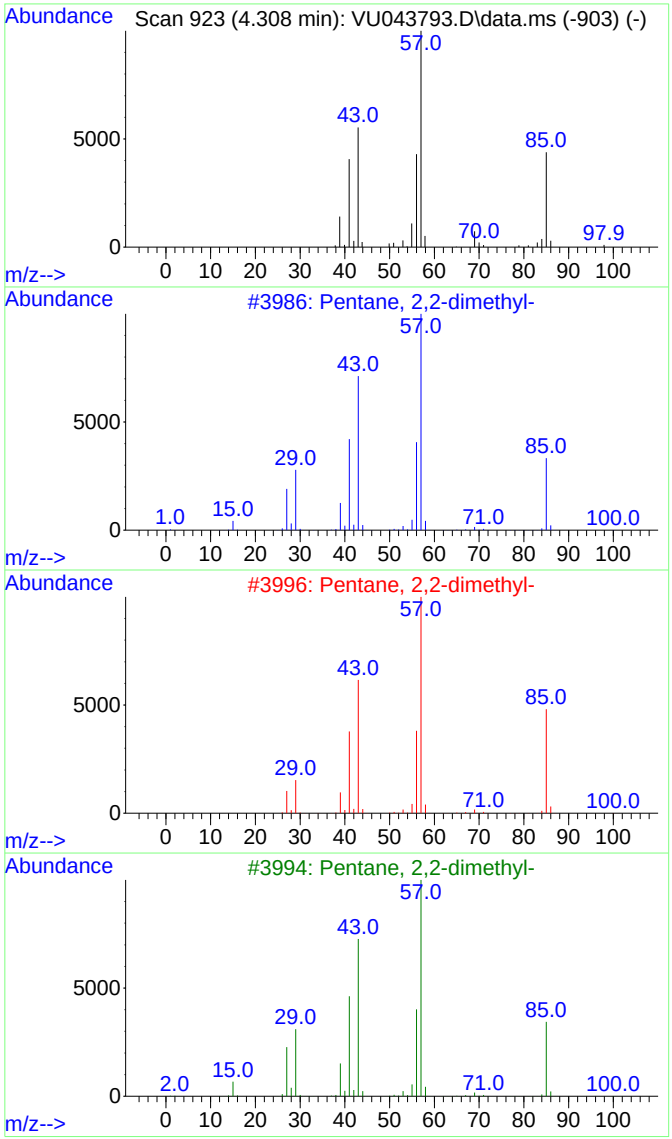
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

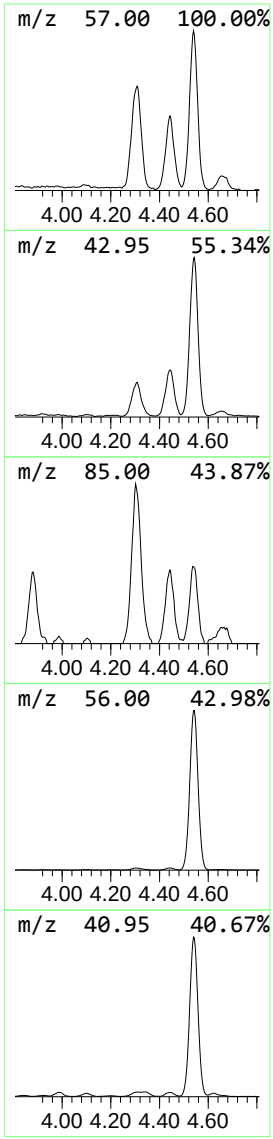
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.308	1.70 ug/L	181095	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	78
2	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	78
3	Pentane, 2,2-dimethyl-	100	C7H16	000590-35-2	64
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59
5	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

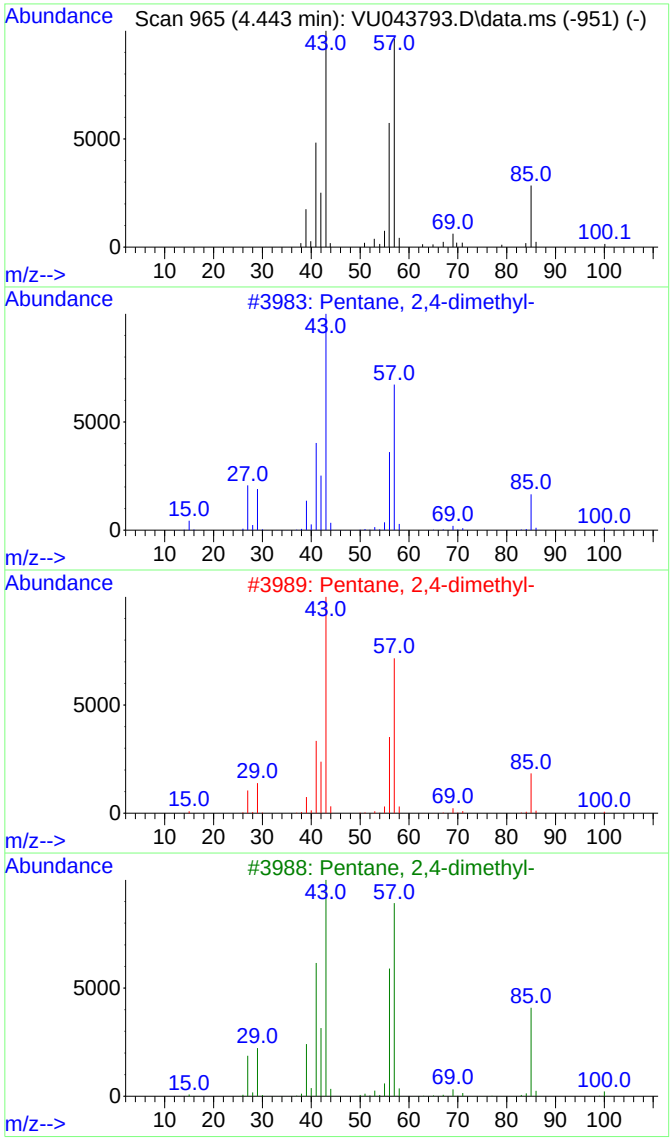
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

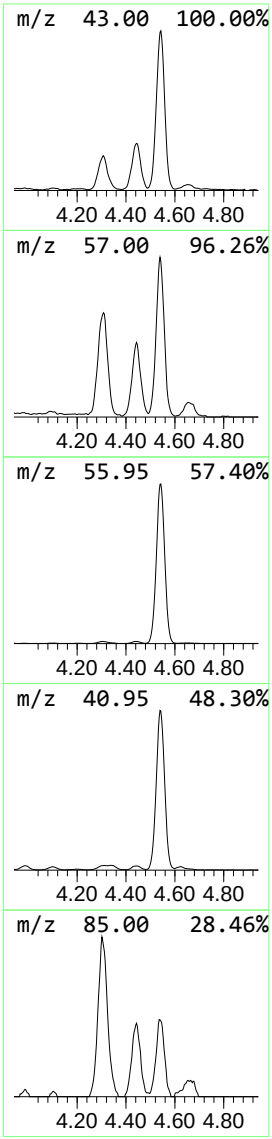
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.443	0.88 ug/L	94206	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	91
2	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	90
3	Pentane, 2,4-dimethyl-	100	C7H16	000108-08-7	72
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	59
5	Pentane, 2,2,3,4-tetramethyl-	128	C9H20	001186-53-4	56





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

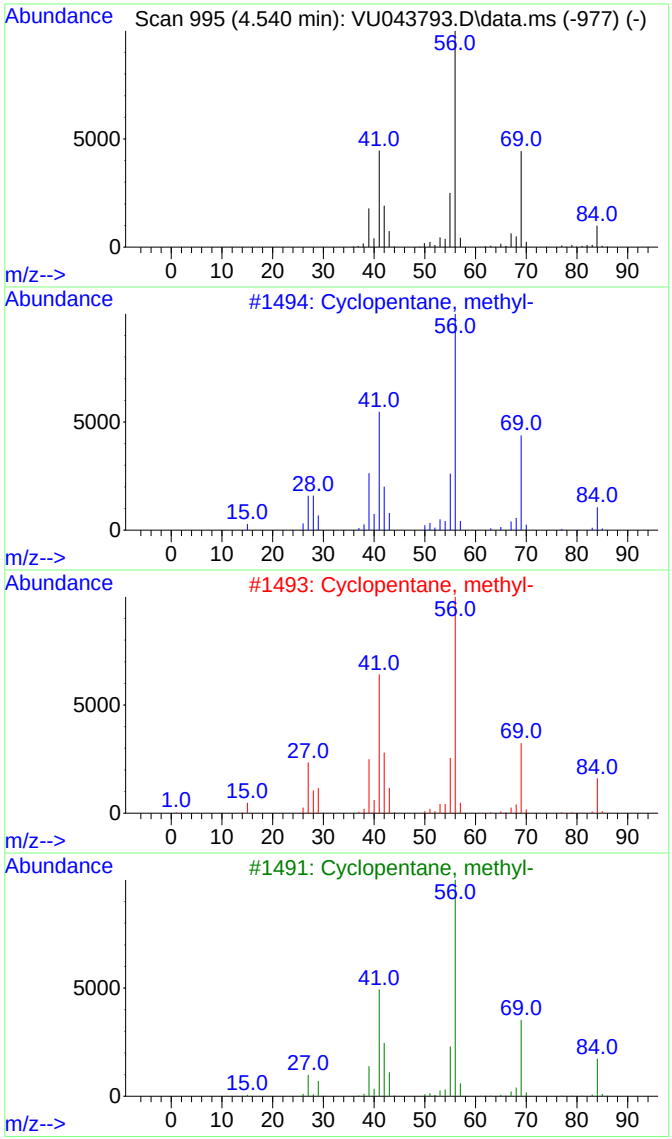
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

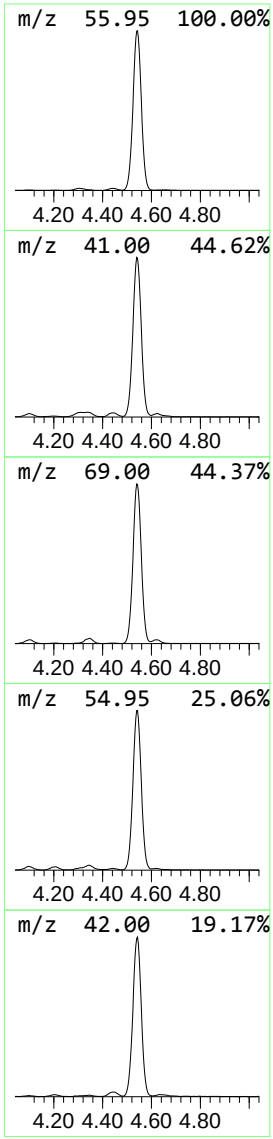
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 (DEL) Alkane: Cyclic4.540 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.540	31.99 ug/L	3417950	1,4-Difluorobenzene	6.256

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	94
2	Cyclopentane, methyl-	84	C6H12	000096-37-7	80
3	Cyclopentane, methyl-	84	C6H12	000096-37-7	78
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

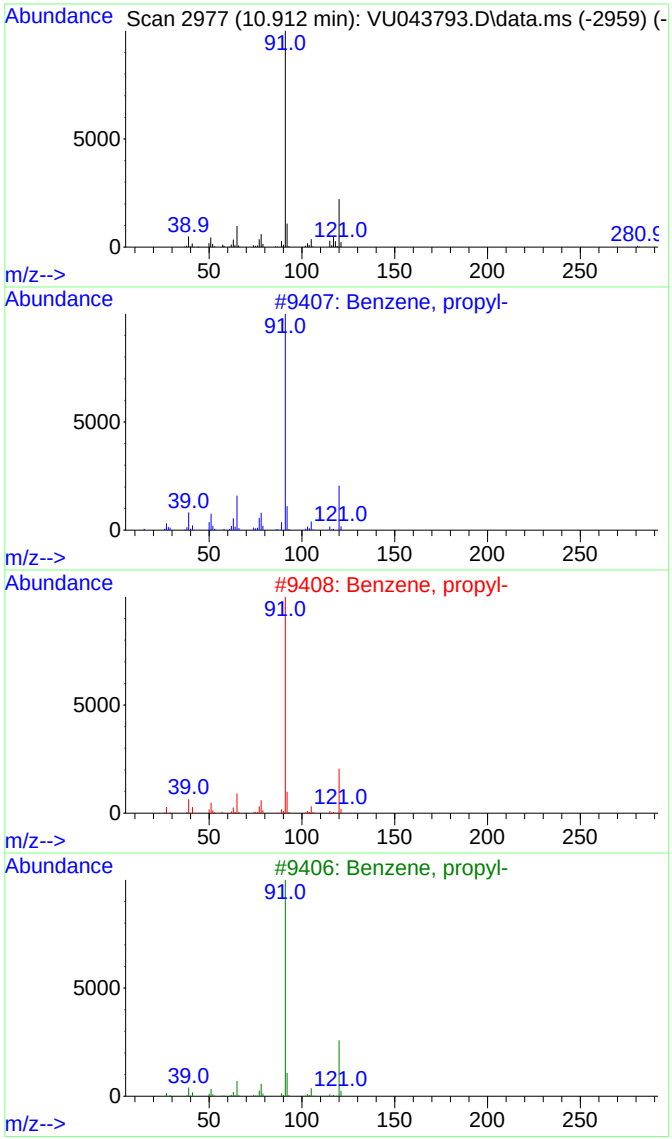
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

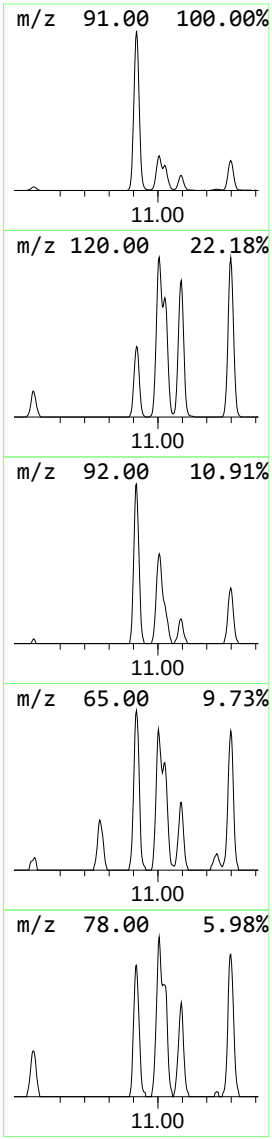
****

Peak Number 8 Benzene, propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.912	1.10 ug/L	192987	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, propyl-	120	C9H12	000103-65-1	87
2	Benzene, propyl-	120	C9H12	000103-65-1	87
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	72
5	Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

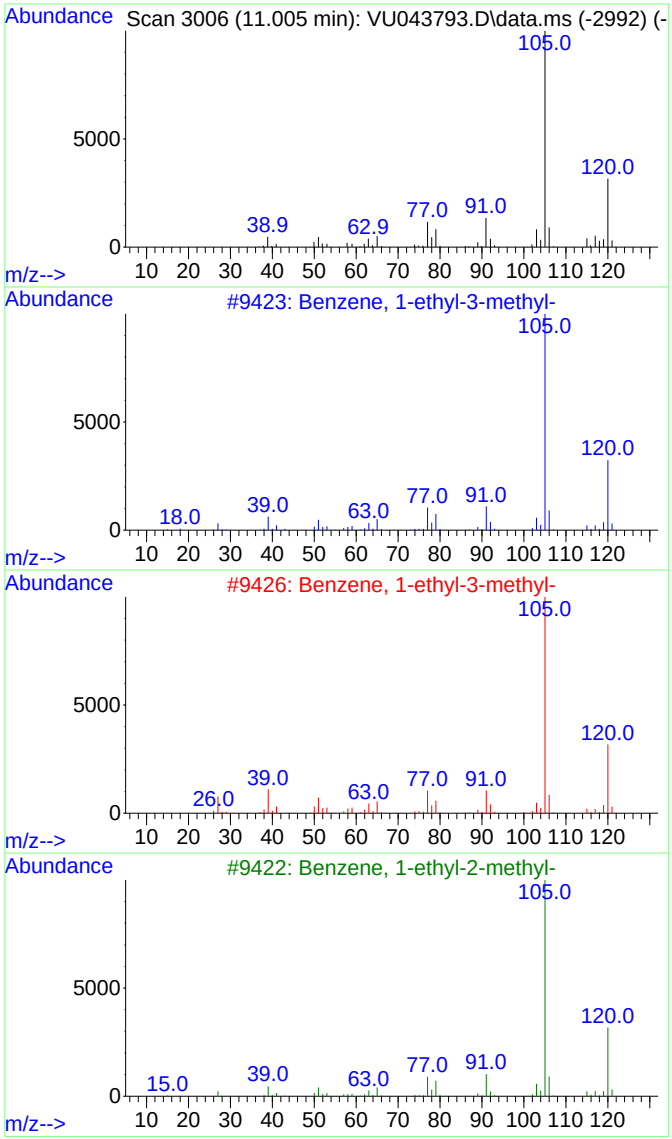
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

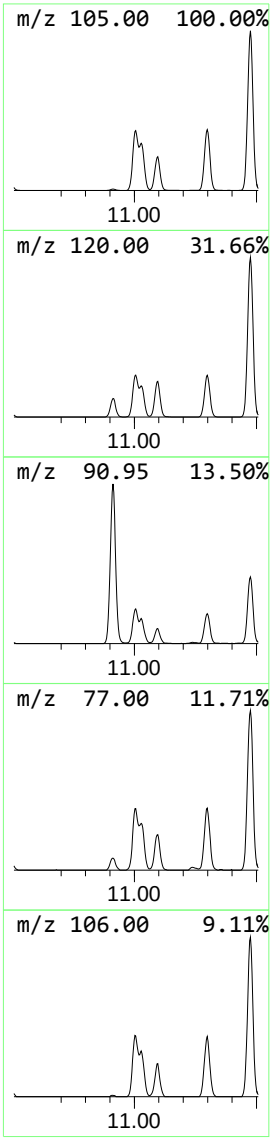
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 9 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.006	2.27 ug/L	396760	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	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-2-methyl-	120	C9H12	000611-14-3	94
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

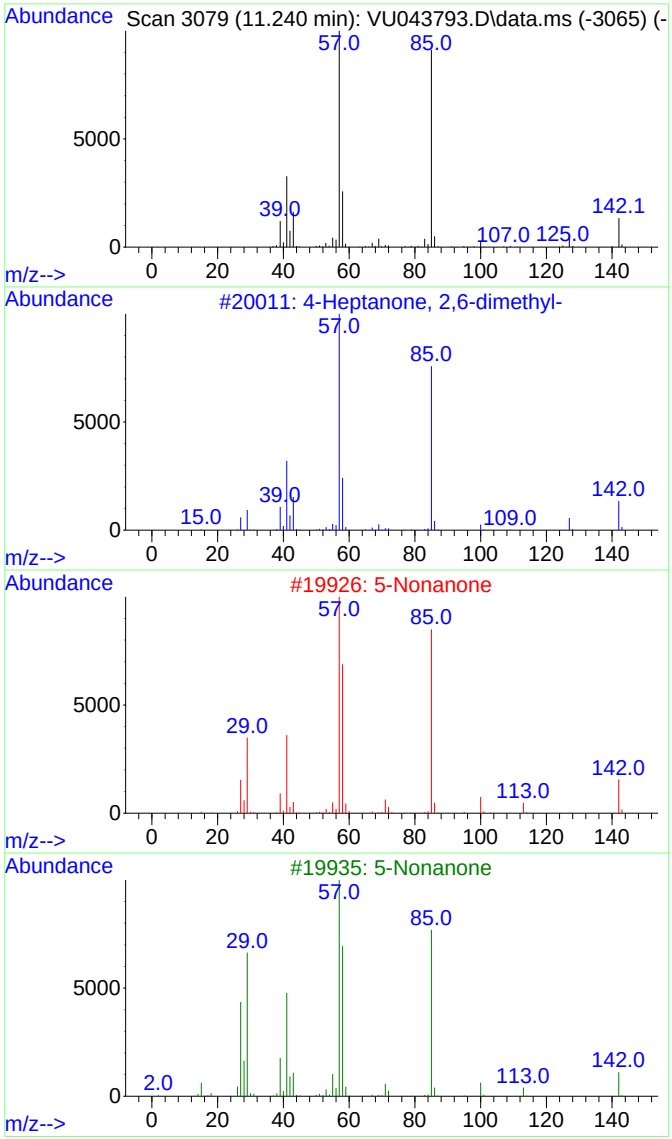
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

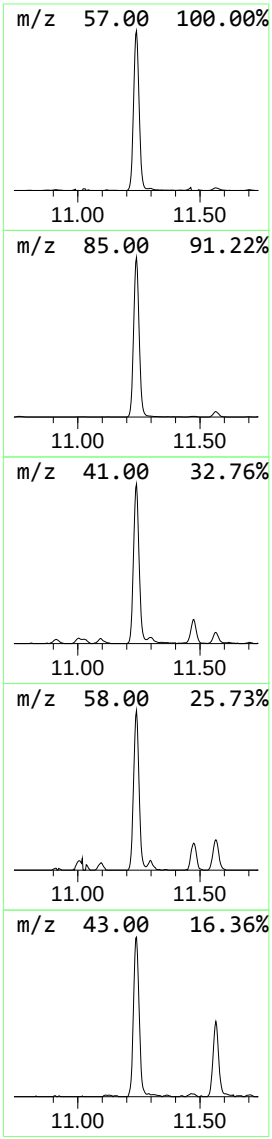
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 4-Heptanone, 2,6-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.240	3.98 ug/L	697272	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	94
2	5-Nonanone	142	C9H18O	000502-56-7	59
3	5-Nonanone	142	C9H18O	000502-56-7	59
4	2,4-Hexanedione, 5,5-dimethyl-	142	C8H14O2	007307-04-2	53
5	4-Heptanone, 2,6-dimethyl-	142	C9H18O	000108-83-8	47





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

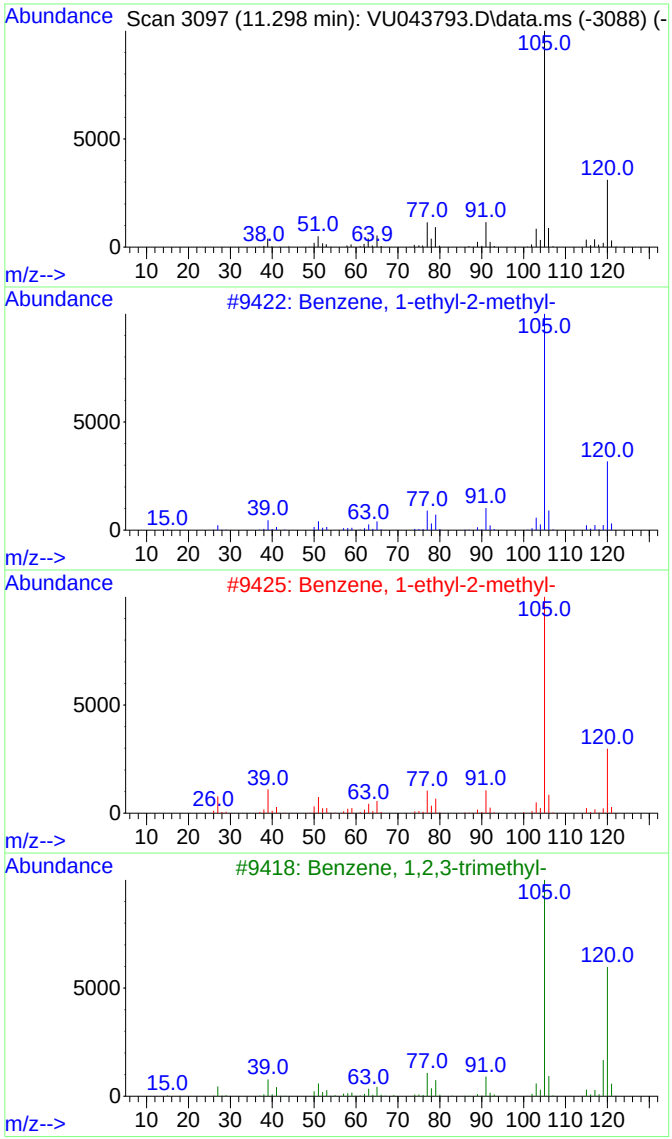
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

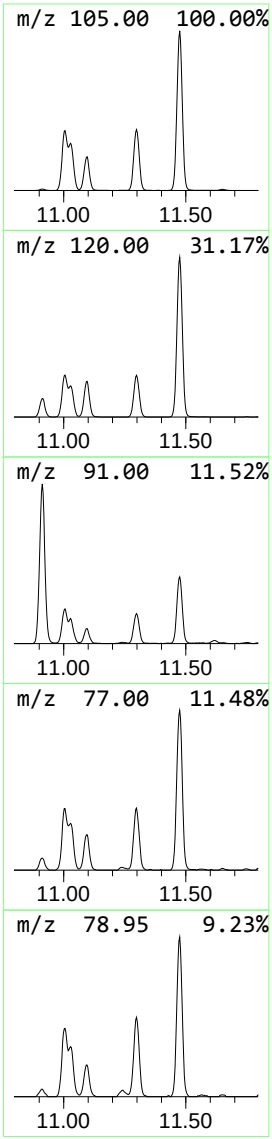
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 1-ethyl-2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.298	2.26 ug/L	395468	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
3	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
4	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

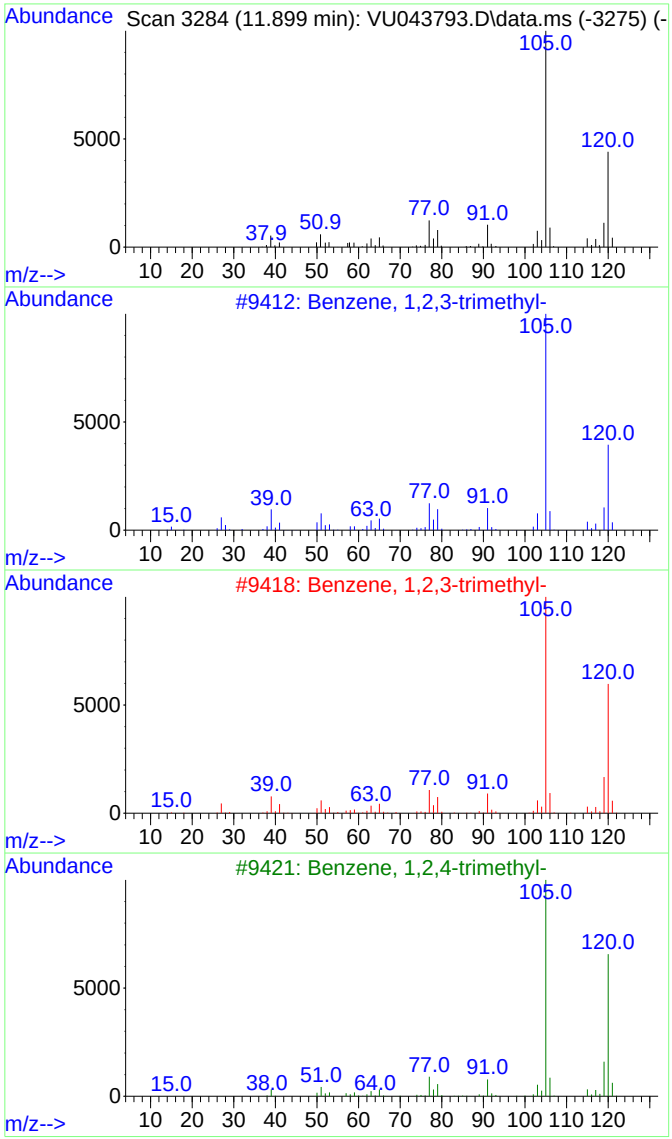
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

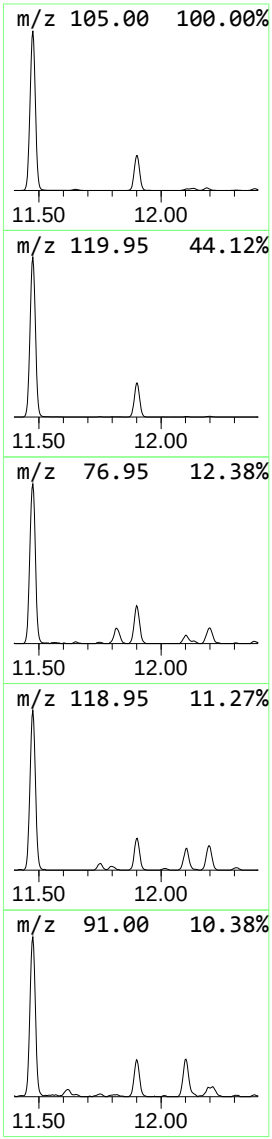
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 12 Benzene, 1,2,3-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.899	1.35 ug/L	236741	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	95
4	Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5	Mesitylene	120	C9H12	000108-67-8	93





Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL-WATER-QT3-2021-06

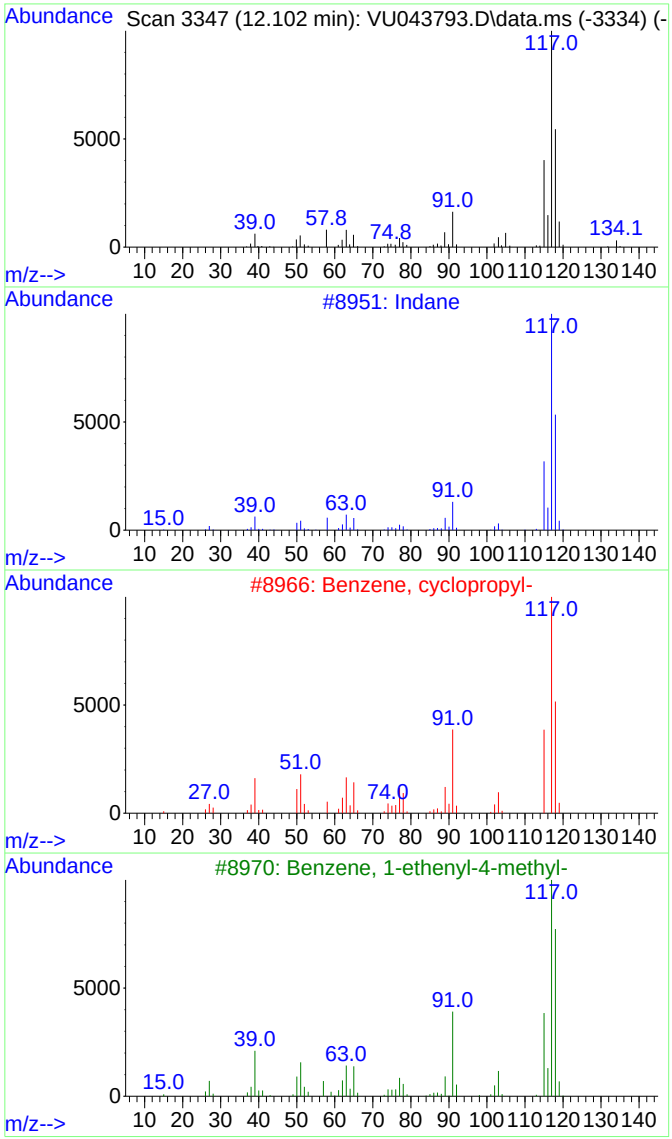
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

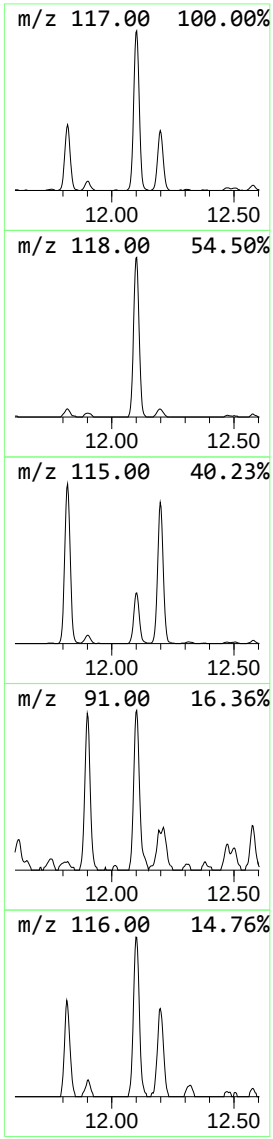
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Indane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.102	1.18 ug/L	206364	1,4-Dichlorobenzene-d4	11.819

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	76
2	Benzene, cyclopropyl-	118	C9H10	000873-49-4	70
3	Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	68
4	trans-Cinnamyl bromide	196	C9H9Br	026146-77-0	64
5	Indane	118	C9H10	000496-11-7	62





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU051721\
Data File : VU043793.D
Acq On : 17 May 2021 15:24
Operator : SY/MD
Sample : M2250-02
Misc : 25.0mL/MSVOA_U/WATER
ALS Vial : 15 Sample Multiplier: 1

Instrument : MSVOA_U
ClientSampleId : MDL-WATER-QT3-2021-06

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR051521WMA.M
Quant Title : TRACE VOA SFAM1.0

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	3.369	27.1	ug/L	2895440	1	6.256	534146	5.0
(DEL) Alkane: S...	3.707	17.3	ug/L	1847490	1	6.256	534146	5.0
(DEL) Alkane: S...	3.986	0.5	ug/L	54858	1	6.256	534146	5.0
(DEL) Alkane: S...	4.308	1.7	ug/L	181095	1	6.256	534146	5.0
(DEL) Alkane: S...	4.443	0.9	ug/L	94206	1	6.256	534146	5.0
(DEL) Alkane: C...	4.540	32.0	ug/L	3417950	1	6.256	534146	5.0
Benzene, propyl-	10.912	1.1	ug/L	192987	3	11.819	875505	5.0
Benzene, 1-ethy...	11.006	2.3	ug/L	396760	3	11.819	875505	5.0
4-Heptanone, 2,...	11.240	4.0	ug/L	697272	3	11.819	875505	5.0
Benzene, 1-ethy...	11.298	2.3	ug/L	395468	3	11.819	875505	5.0
Benzene, 1,2,3-...	11.899	1.4	ug/L	236741	3	11.819	875505	5.0
Indane	12.102	1.2	ug/L	206364	3	11.819	875505	5.0