LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

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\SFAMVTR112321WMA.M
Title : TRACE VOA SFAM1.0

Signal : TIC: VV023819.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.307	76	83	94	rVB	61232	60859	7.66%	0.812%
2	1.568	157	164	175	rBV	52825	59790	7.53%	0.798%
3	1.632	175	184	200	rVB	48971	62795	7.91%	0.838%
4	1.806	229	238	251	rVB	58491	75447	9.50%	1.006%
5	2.108	316	332	350	rBV	103694	159220	20.05%	2.124%
6	2.436	424	434	444	rBV7	5132	11381	1.43%	0.152%
7	2.529	444	463	473	rVV3	38107	116155	14.62%	1.549%
8	2.581	474	479	498	rVB8	9252	19687	2.48%	0.263%
9	2.764	520	536	550	rBV2	28064	57334	7.22%	0.765%
10	2.953	578	595	608	rBV4	10926	23843	3.00%	0.318%
11	3.040	610	622	640	rVB2	28299	57146	7.19%	0.762%
12	3.764	830	847	865	rBV2	48193	122389	15.41%	1.632%
13	3.912	880	893	923	rBV4	38569	138991	17.50%	1.854%
14	4.346	1013	1028	1064	rVB	78547	224947	28.32%	3.000%
15	4.677	1115	1131	1146	rBV4	43272	107050	13.48%	1.428%
16	4.912	1191	1204	1208	rBV6	4774	9557	1.20%	0.127%
17	5.047	1229	1246	1273	rBV2	171017	433555	54.58%	5.783%
18	5.317	1318	1330	1344	rVB5	9174	19994	2.52%	0.267%
19	5.619	1412	1424	1453	rBV	143204	304270	38.31%	4.058%
20	5.915	1504	1516	1543	rBV	119018	245504	30.91%	3.275%
21	6.069	1550	1564	1577	rBV	95199	202970	25.55%	2.707%
22	6.130	1579	1583	1600	rVB3	20871	38019	4.79%	0.507%
23	6.989	1839	1850	1872	rVB2	47452	88121	11.09%	1.175%
24	7.317	1941	1952	1971	rBV	216137	373947	47.08%	4.988%
25	7.625	2038	2048	2073	rBV2	28277	53536	6.74%	0.714%
26	7.976	2138	2157	2182	rBV	479137	794288	100.00%	10.595%
27	8.092	2182	2193	2230	rBV	155047	363729	45.79%	4.852%
28	8.854	2418	2430	2454	rBV	213875	355359	44.74%	4.740%
29	9.014	2471	2480	2493	rBV	33292	52168	6.57%	0.696%
30	9.136	2505	2518	2535	rBV	150556	252912	31.84%	3.373%
31	9.545	2635	2645	2660	rBV2	20399	32859	4.14%	0.438%
32	9.937	2759	2767	2777	rBV2	6804	10774	1.36%	0.144%
33	10.217	2843	2854	2870	rBV	69983	113684	14.31%	1.516%
34	10.355	2883	2897	2915	rBV	24924	43081	5.42%	0.575%
35	10.448	2915	2926	2945	rVV	117911	263625	33.19%	3.516%
36	10.538	2945	2954	2972	rVB	68396	105333	13.26%	1.405%

LSC Area Percent Report

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

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\SFAMVTR112321WMA.M
Title : TRACE VOA SFAM1.0

37	10.738	3006	3016	3034	rBV	69057	109247	13.75%	1.457%
38	10.915	3058	3071	3088	rBV	334774	506846	63.81%	6.761%
39	11.249	3163	3175	3192	rBV	248870	386481	48.66%	5.155%
40	11.336	3192	3202	3216	rVB	71927	111416	14.03%	1.486%
41	11.529	3251	3262	3271	rBV	33141	60372	7.60%	0.805%
42	11.574	3272	3276	3282	rVV	14036	18795	2.37%	0.251%
43	11.625	3282	3292	3315	rVB2	243295	414997	52.25%	5.535%
44	11.821	3344	3353	3364	rBV2	5422	8059	1.01%	0.107%
45	11.911	3371	3381	3386	rBV3	18018	27937	3.52%	0.373%
46	11.940	3386	3390	3400	rVB2	15619	21532	2.71%	0.287%
47	12.017	3405	3414	3421	rBV2	29544	47373	5.96%	0.632%
48	12.114	3435	3444	3459	rBV2	37867	61749	7.77%	0.824%
49	12.316	3495	3507	3516	rBV4	7163	10833	1.36%	0.144%
50	12.413	3527	3537	3545	rBV2	19728	28758	3.62%	0.384%
51	12.464	3545	3553	3564	rVB	26373	39683	5.00%	0.529%
52	12.725	3624	3634	3648	rVB2	24471	38092	4.80%	0.508%
53	12.882	3665	3683	3695	rBV3	46592	78355	9.86%	1.045%
54	13.050	3727	3735	3745	rBV9	5379	9608	1.21%	0.128%
55	13.217	3778	3787	3795	rVB3	7199	11893	1.50%	0.159%
56	13.271	3795	3804	3810	rBV5	7167	10375	1.31%	0.138%
57	13.506	3867	3877	3886	rBV	18371	28806	3.63%	0.384%
58	14.638	4217	4229	4239	rBV	7175	11586	1.46%	0.155%

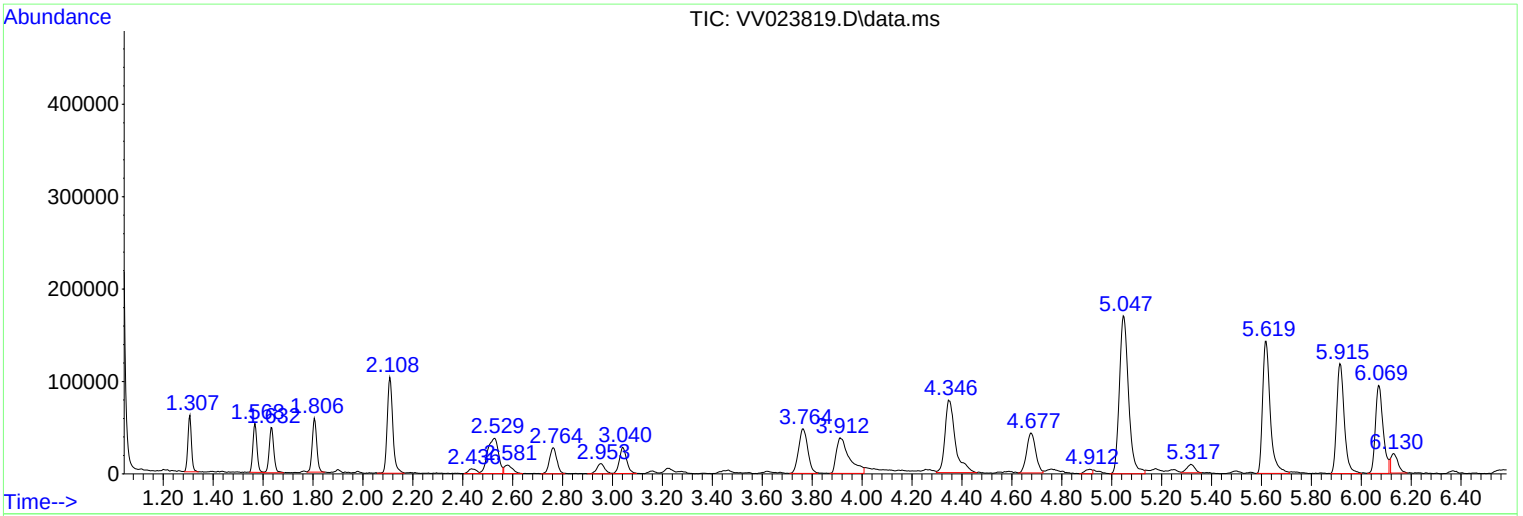
Sum of corrected areas: 7497112

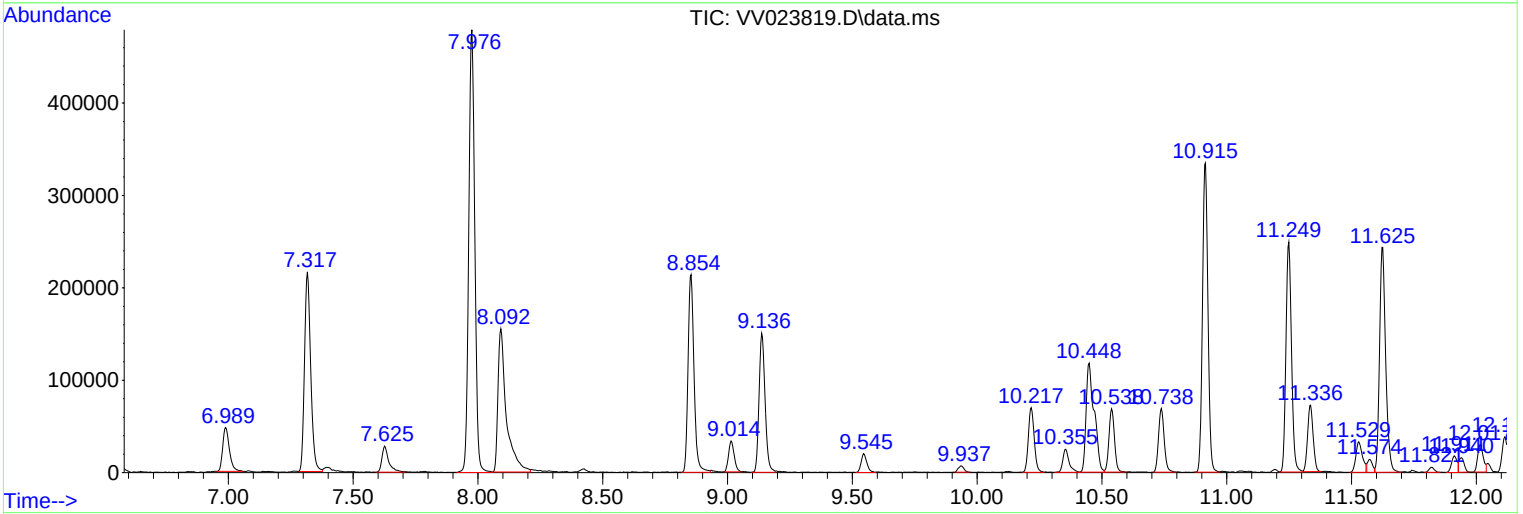
Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

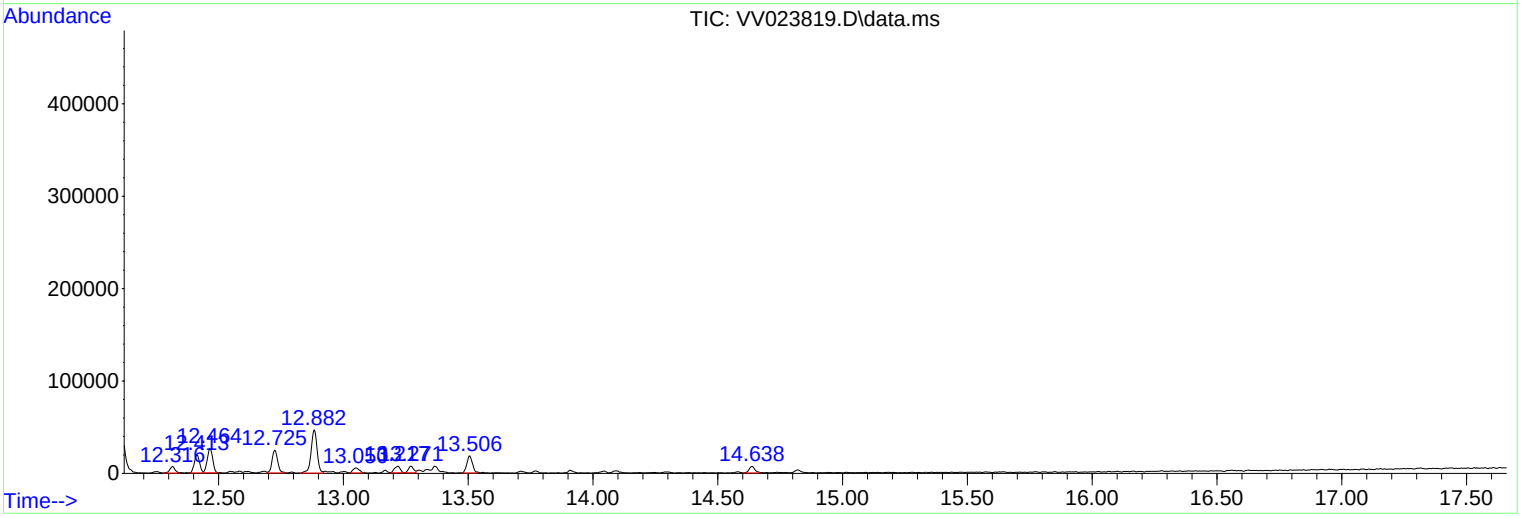
Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.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\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

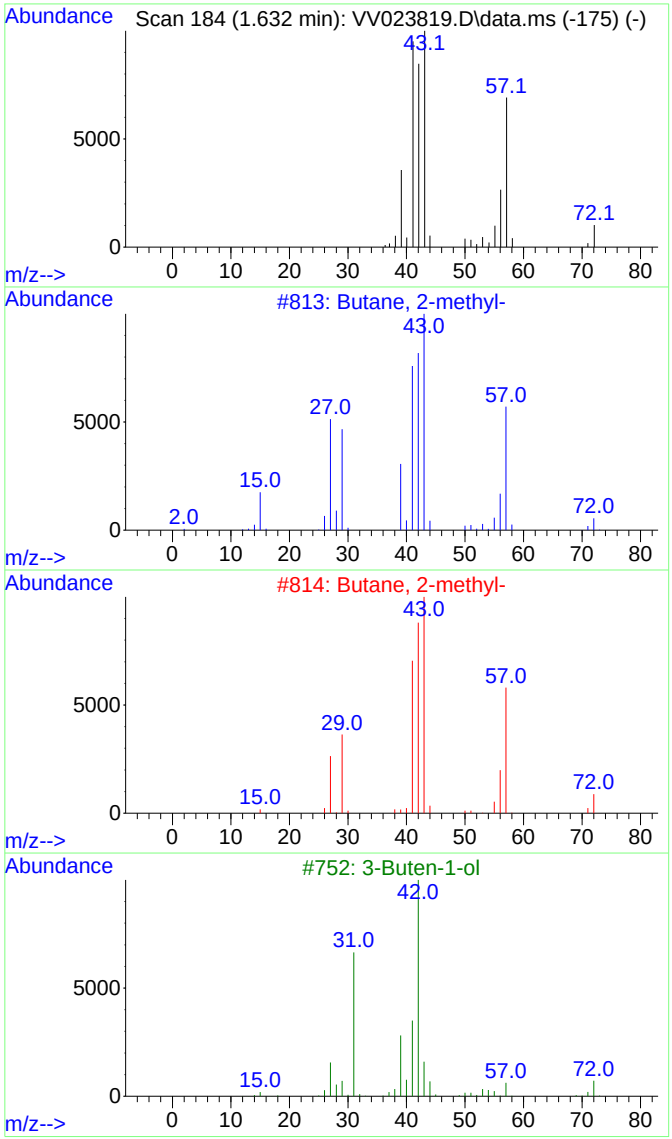
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

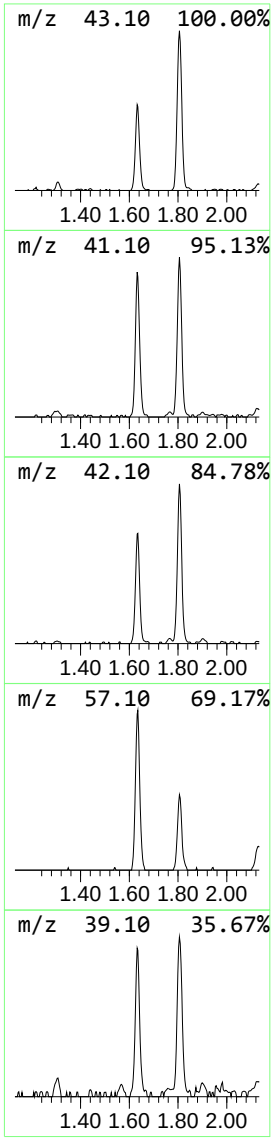
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.632	1.03 ug/L	62795	1,4-Difluorobenzene	5.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Butane, 2-methyl-	72	C5H12	000078-78-4	80
2	Butane, 2-methyl-	72	C5H12	000078-78-4	80
3	3-Buten-1-ol	72	C4H8O	000627-27-0	43
4	1,3-Dimethyldiaziridine	72	C3H8N2	026177-36-6	33
5	Pentane	72	C5H12	000109-66-0	28





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

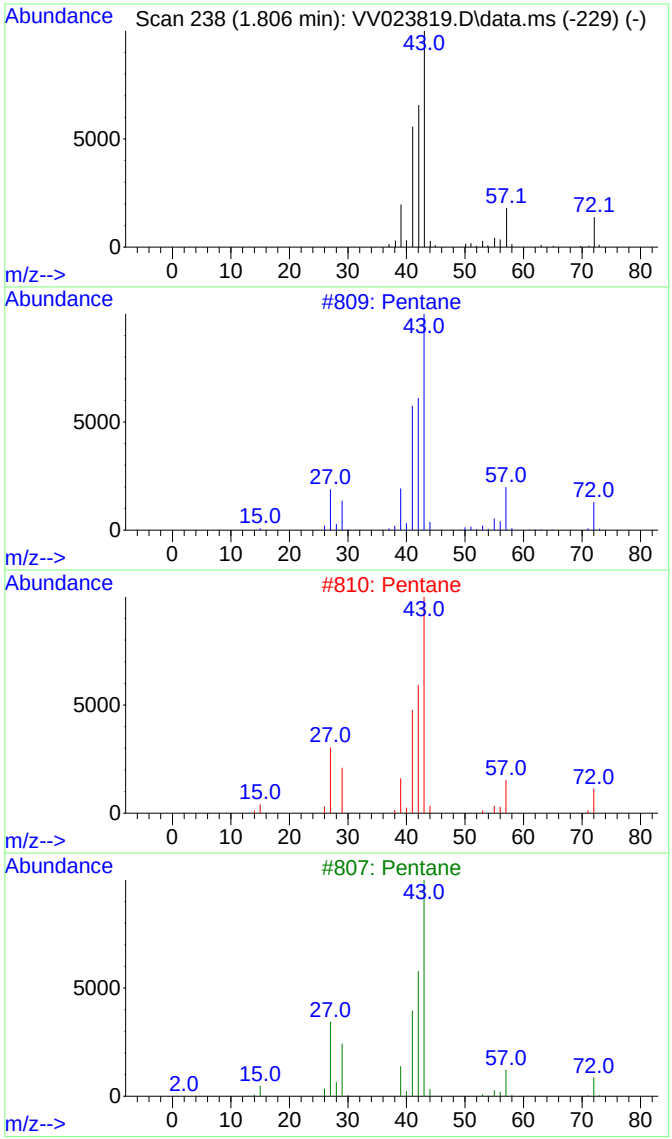
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

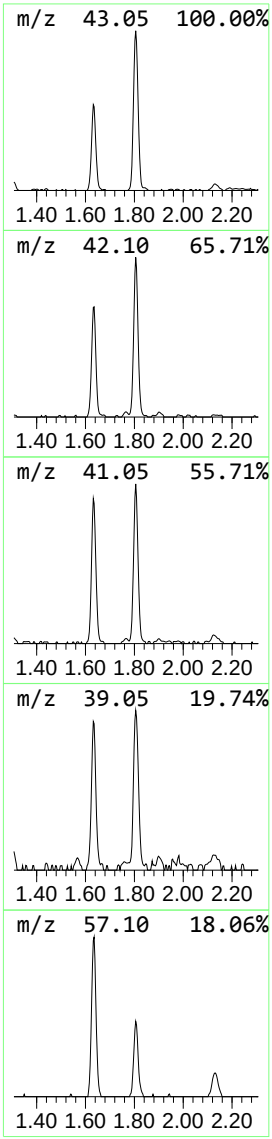
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
1.806	1.24 ug/L	75447	1,4-Difluorobenzene	5.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane	72	C5H12	000109-66-0	91
2	Pentane	72	C5H12	000109-66-0	86
3	Pentane	72	C5H12	000109-66-0	86
4	Pentane	72	C5H12	000109-66-0	80
5	Butane, 2-methyl-	72	C5H12	000078-78-4	64





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

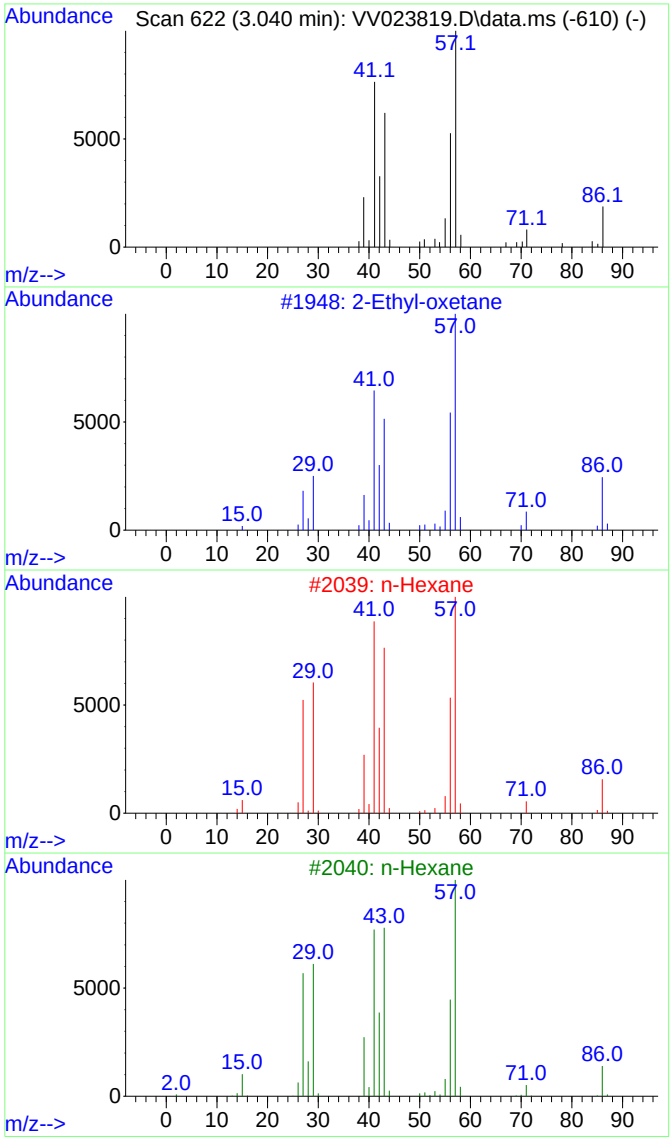
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

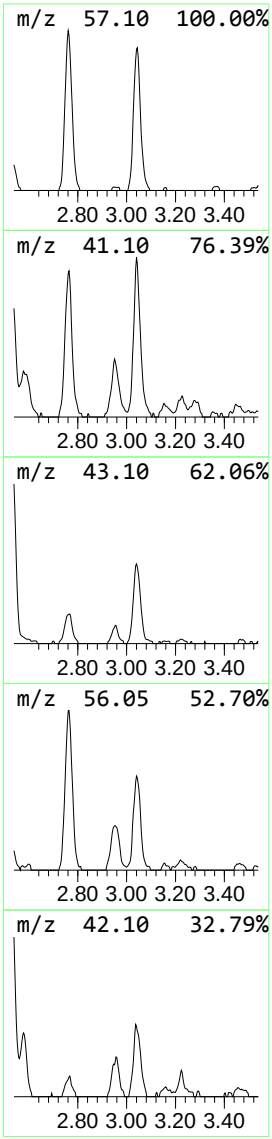
****

Peak Number 3 2-Ethyl-oxetane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.040	0.94 ug/L	57146	1,4-Difluorobenzene	5.616

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Ethyl-oxetane	86	C5H10O	1010386-40-2	91
2	n-Hexane	86	C6H14	000110-54-3	91
3	n-Hexane	86	C6H14	000110-54-3	72
4	n-Hexane	86	C6H14	000110-54-3	64
5	n-Hexane	86	C6H14	000110-54-3	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

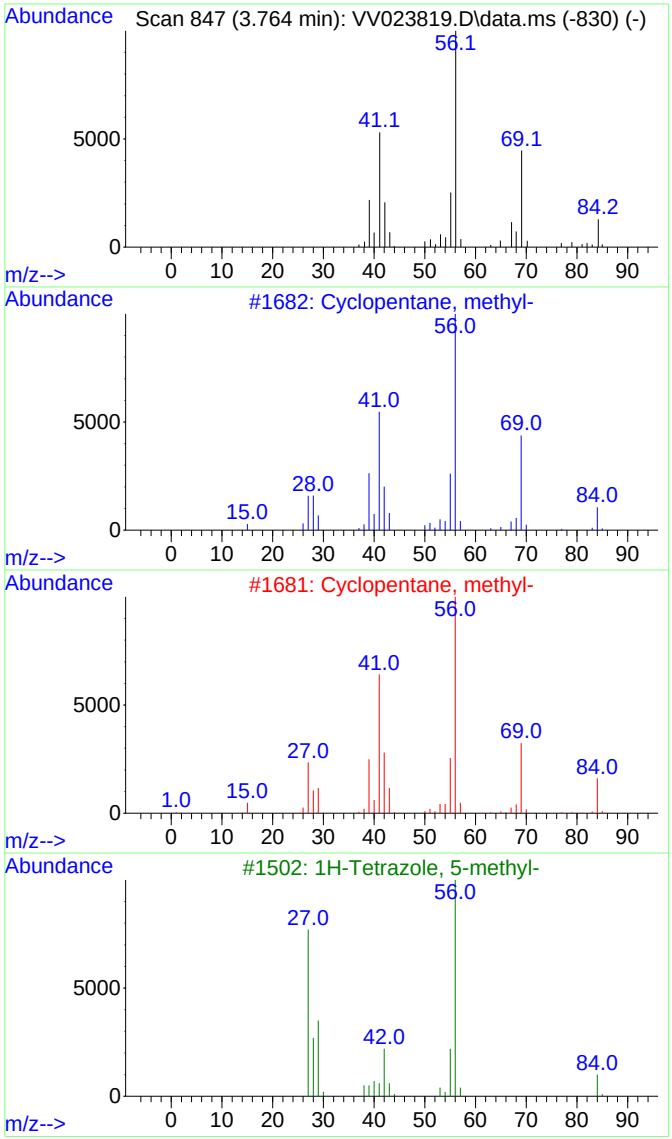
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

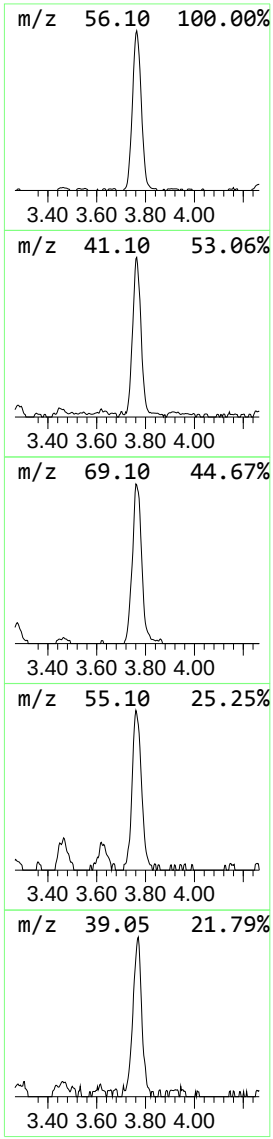
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 (DEL) Alkane: Cyclic3.764 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.764	2.01 ug/L	122389	1,4-Difluorobenzene	5.616

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	80
3	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	72
4	Cyclopentane, methyl-	84	C6H12	000096-37-7	64
5	Cyclobutane, ethyl-	84	C6H12	004806-61-5	59





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

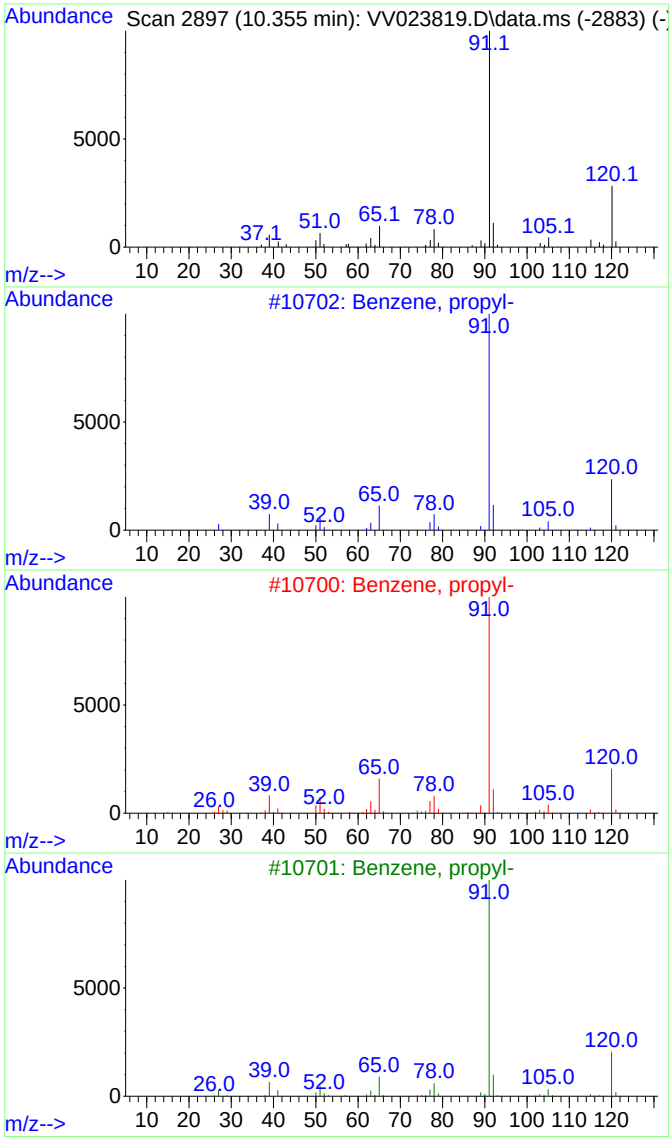
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

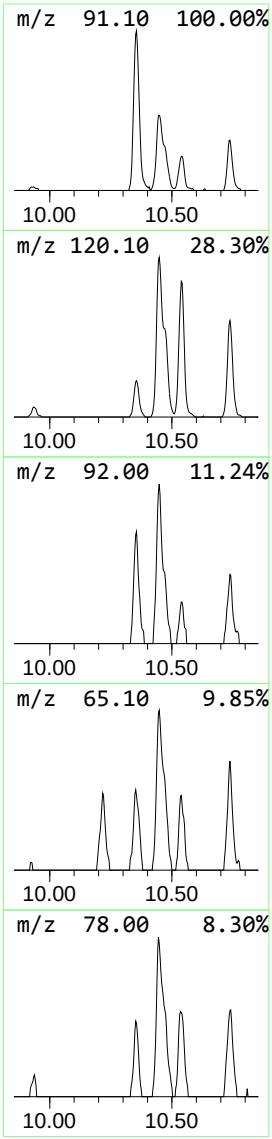
****

Peak Number 6 Benzene, propyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.355	0.56 ug/L	43081	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	93
3	Benzene, propyl-	120	C9H12	000103-65-1	87
4	Benzene, propyl-	120	C9H12	000103-65-1	83
5	N-Benzyl-N-hexylamine	191	C13H21N	025468-44-4	72





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

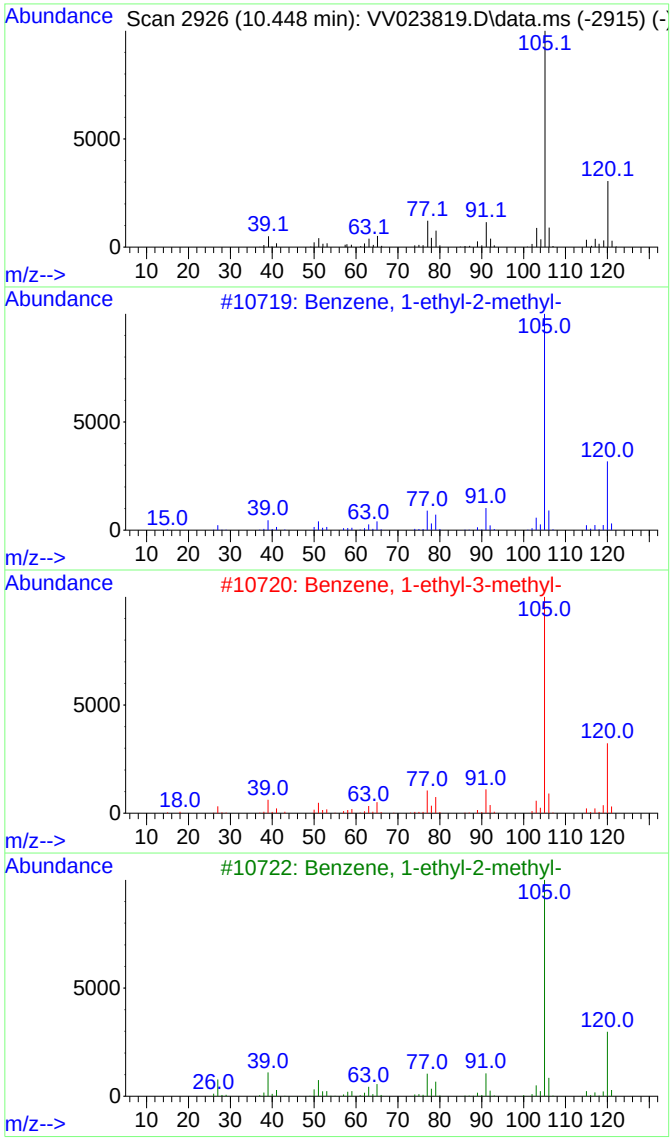
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

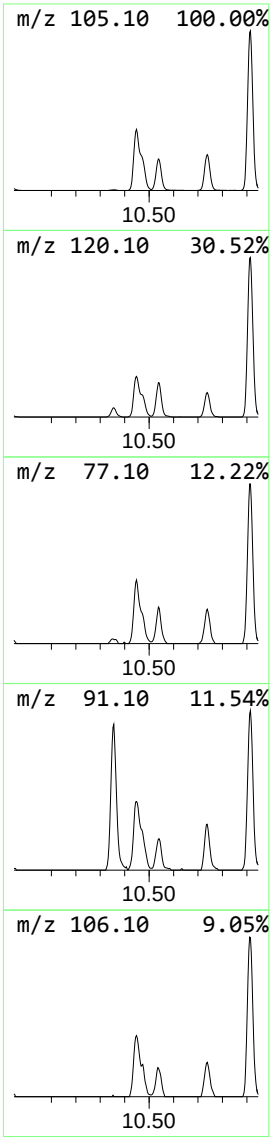
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.448	3.41 ug/L	263625	1,4-Dichlorobenzene-d4	11.249

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	95
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-3-methyl-	120	C9H12	000620-14-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

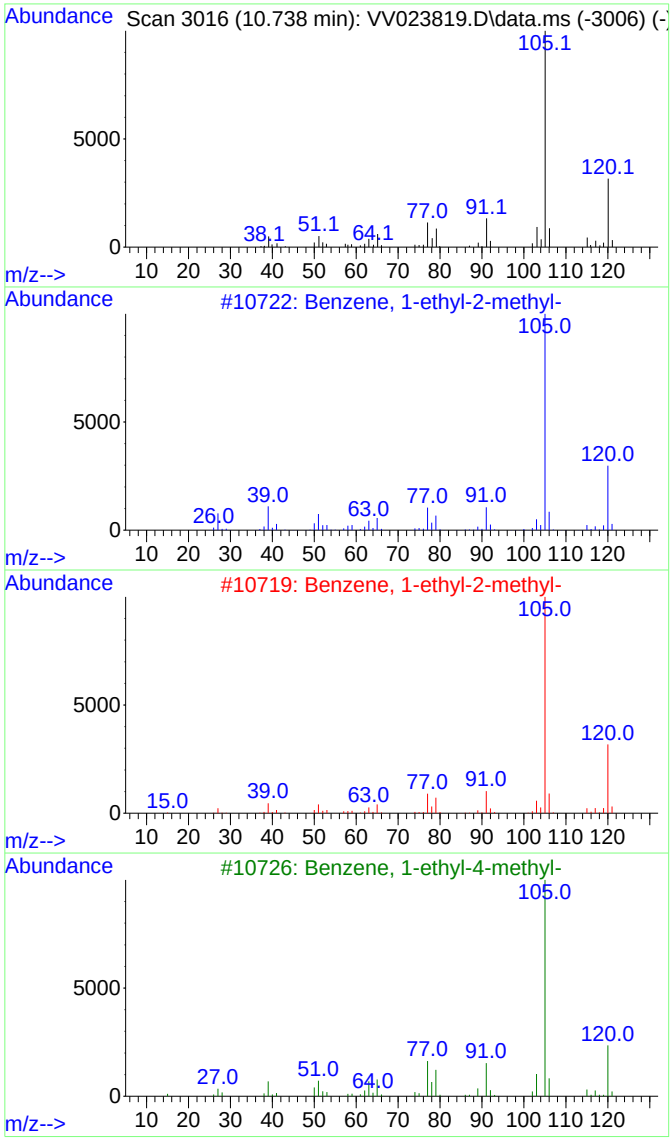
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

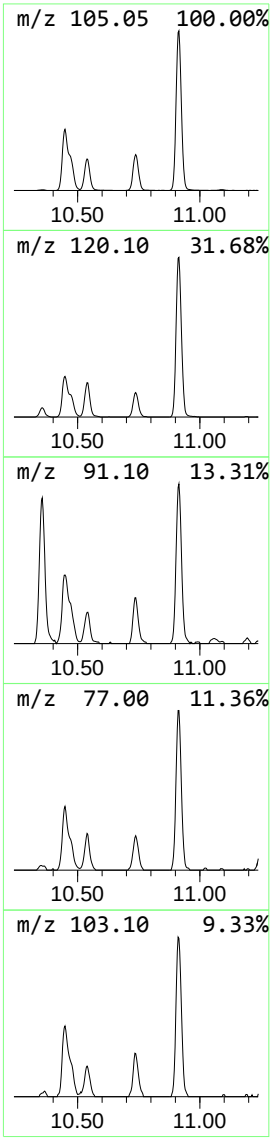
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 Benzene, 1-ethyl-3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.738	1.41 ug/L	109247	1,4-Dichlorobenzene-d4	11.249

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-2-methyl-	120	C9H12	000611-14-3	95
3	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Mesitylene	120	C9H12	000108-67-8	91





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Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

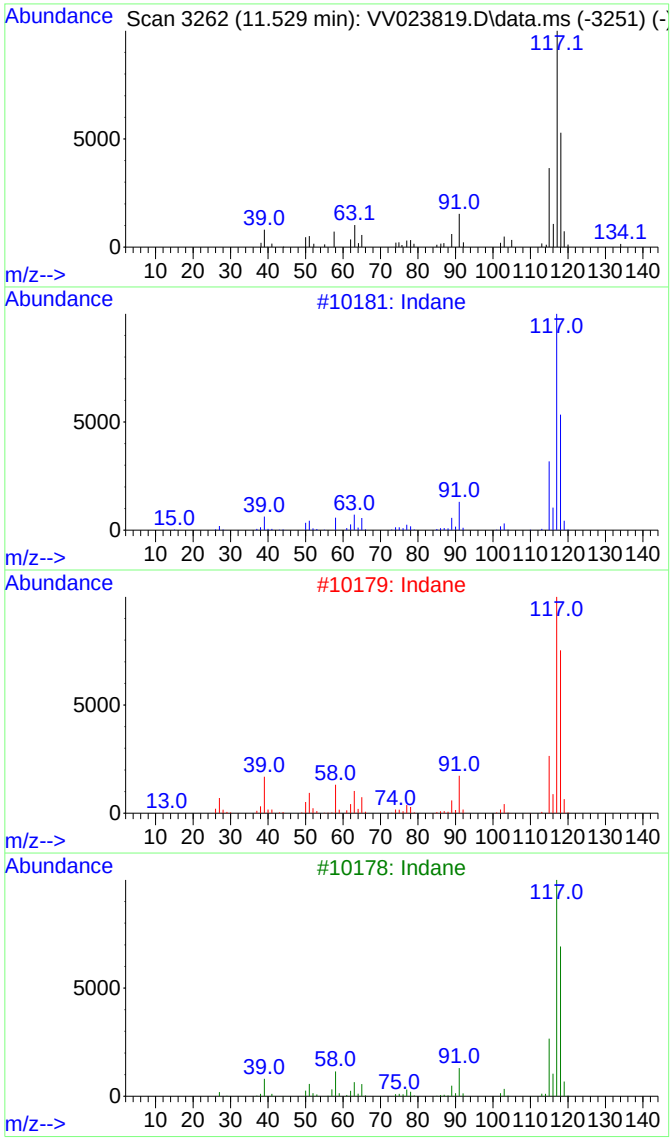
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

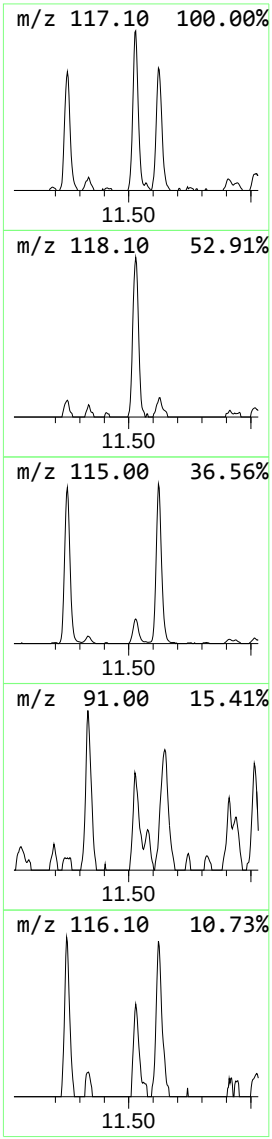
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 Indane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	0.78 ug/L	60372	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	93
2	Indane	118	C9H10	000496-11-7	87
3	Indane	118	C9H10	000496-11-7	81
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
5	Indane	118	C9H10	000496-11-7	81





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

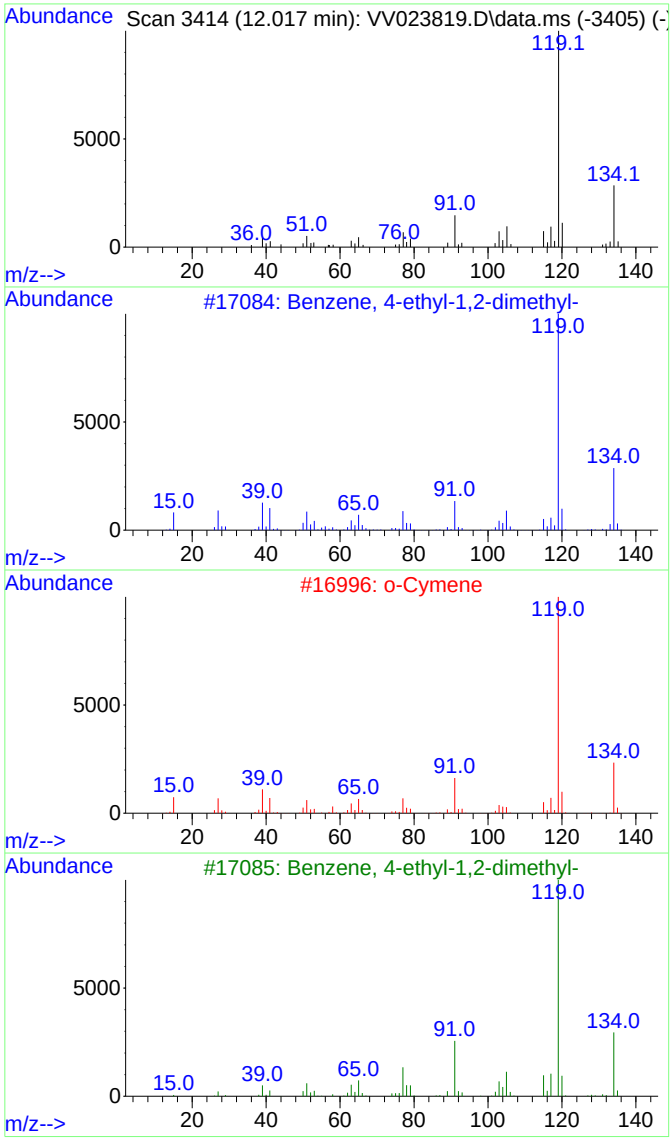
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

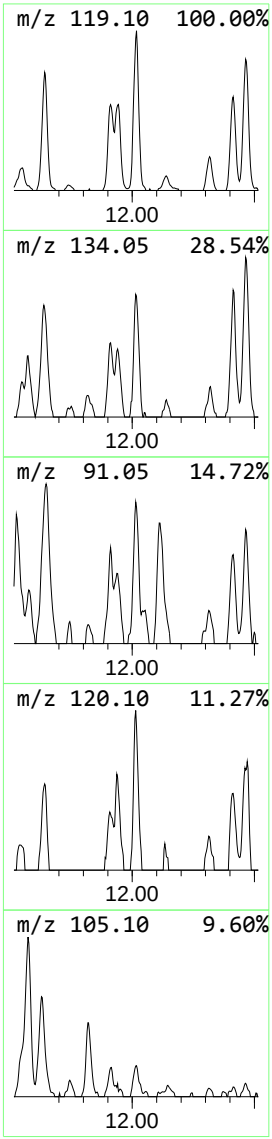
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 11 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.017	0.61 ug/L	47373	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
5	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

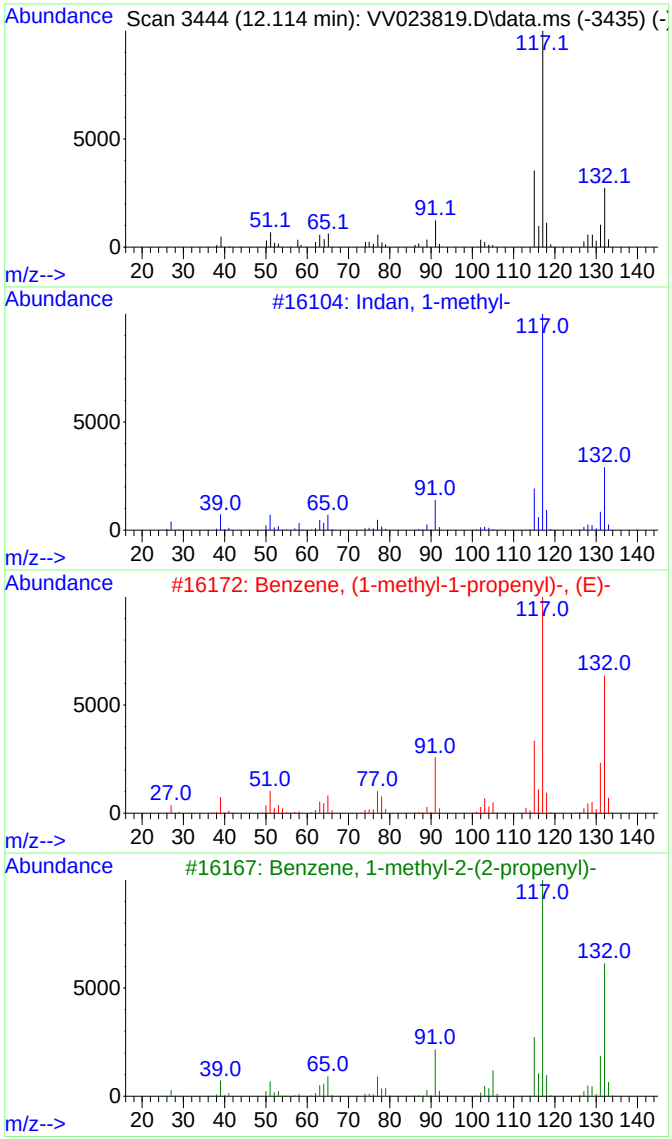
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

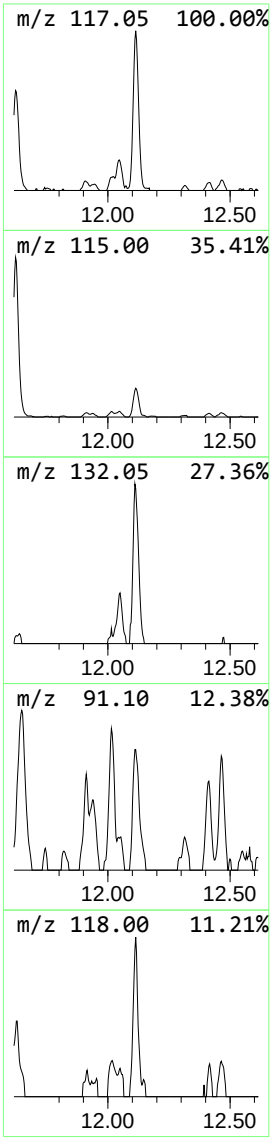
****

Peak Number 12 Indan, 1-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.114	0.80 ug/L	61749	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indan, 1-methyl-	132	C10H12	000767-58-8	90
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	90
3	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	87
4	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
5	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

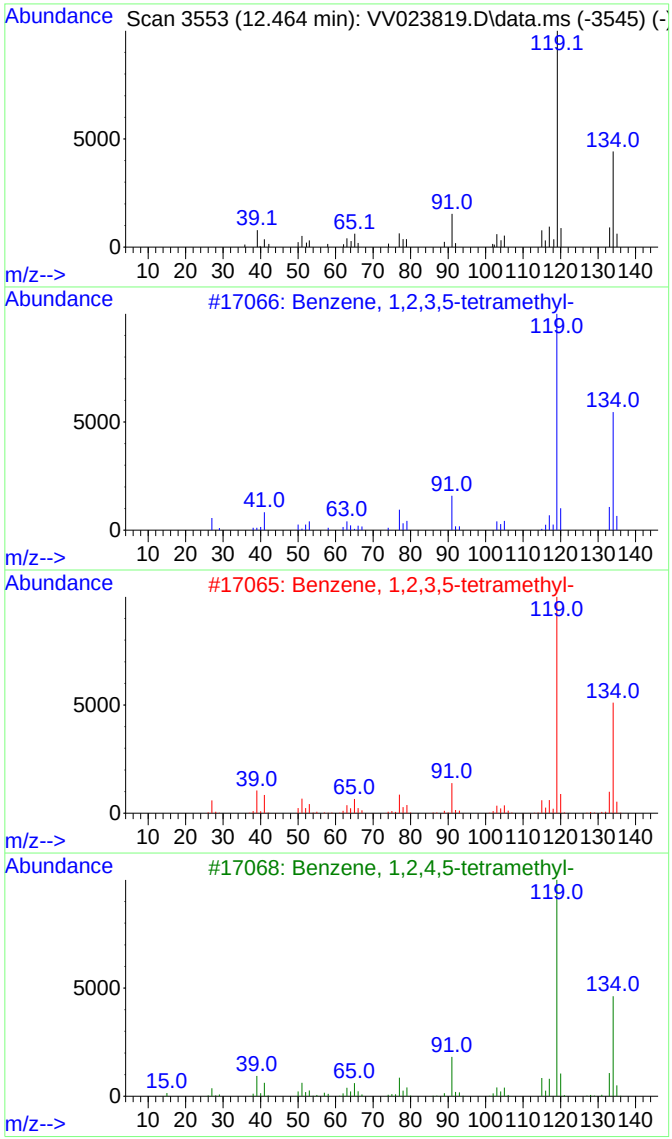
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

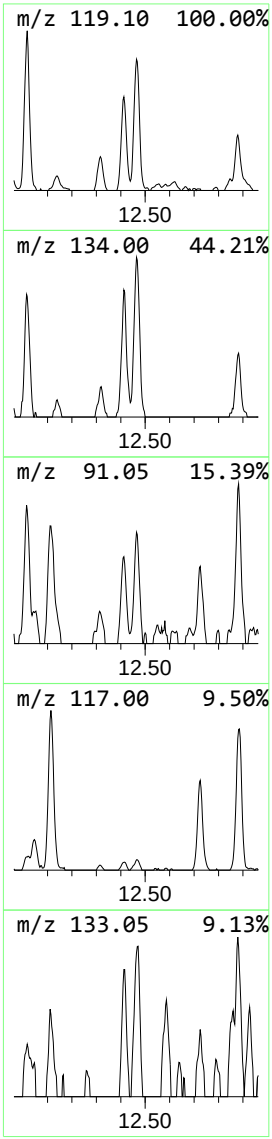
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 13 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.464	0.51 ug/L	39683	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	96
2	Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5	Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95





Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

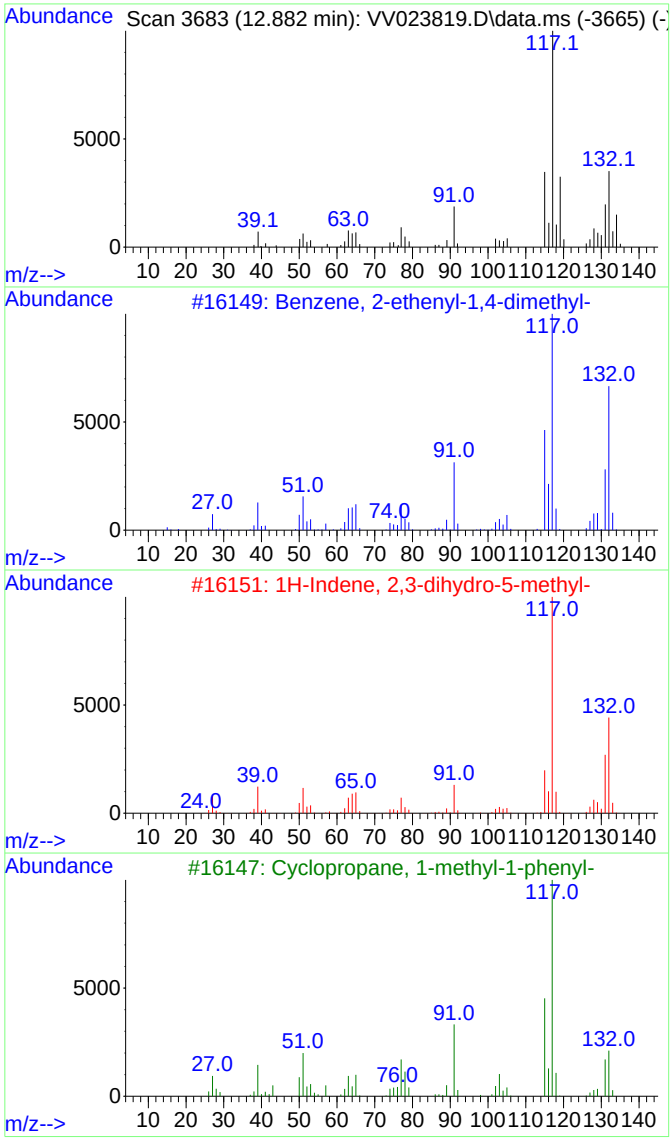
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.M
Quant Title : TRACE VOA SFAM1.0

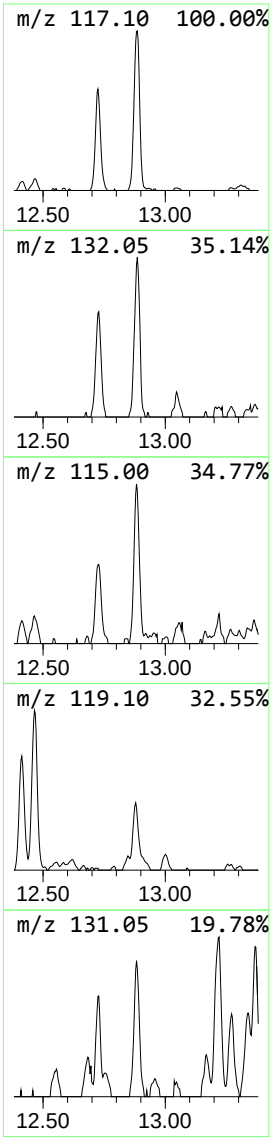
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 14 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.882	1.01 ug/L	78355	1,4-Dichlorobenzene-d4	11.249

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	76
3	Cyclopropane, 1-methyl-1-phenyl-	132	C10H12	002214-14-4	74
4	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	70
5	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	68





Tentatively Identified Compound (LSC) summary

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV120721\
Data File : VV023819.D
Acq On : 07 Dec 2021 13:44
Operator : SY/MD
Sample : M4879-07DL 200X
Misc : 25.0mL/MSVOA_V/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_V
ClientSampleId :
C0G35DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR112321WMA.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
(DEL) Alkane: S...	1.632	1.0	ug/L	62795	1	5.616	304270	5.0
(DEL) Alkane: S...	1.806	1.2	ug/L	75447	1	5.616	304270	5.0
2-Ethyl-oxetane	3.040	0.9	ug/L	57146	1	5.616	304270	5.0
(DEL) Alkane: C...	3.764	2.0	ug/L	122389	1	5.616	304270	5.0
Benzene, propyl-	10.355	0.6	ug/L	43081	3	11.249	386481	5.0
Benzene, 1-ethy...	10.448	3.4	ug/L	263625	3	11.249	386481	5.0
Benzene, 1-ethy...	10.738	1.4	ug/L	109247	3	11.249	386481	5.0
Indane	11.529	0.8	ug/L	60372	3	11.249	386481	5.0
Benzene, 4-ethy...	12.017	0.6	ug/L	47373	3	11.249	386481	5.0
Indan, 1-methyl-	12.114	0.8	ug/L	61749	3	11.249	386481	5.0
Benzene, 1,2,3,...	12.464	0.5	ug/L	39683	3	11.249	386481	5.0
Benzene, 2-ethe...	12.882	1.0	ug/L	78355	3	11.249	386481	5.0