

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VV032522\
 Data File : VV025178.D
 Acq On : 25 Mar 2022 14:55
 Operator : SY/MD
 Sample : N2083-07DL 10X
 Misc : 25mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 EW607DL

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\SFAMVTR031522WMA.M
 Title : TRACE VOA SFAM1.0

Signal : TIC: VV025178.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	98	rVB	65069	67200	14.32%	1.608%
2	1.587	157	170	186	rVB2	83497	154606	32.95%	3.700%
3	2.108	322	332	350	rBV	96962	142035	30.27%	3.400%
4	2.510	443	457	469	rBV2	18666	31605	6.74%	0.756%
5	3.191	655	669	691	rBV	51288	108103	23.04%	2.587%
6	3.902	880	890	928	rBV	49632	182456	38.89%	4.367%
7	4.352	1009	1030	1057	rBV2	85228	217401	46.34%	5.203%
8	4.613	1095	1111	1128	rBV2	29361	73705	15.71%	1.764%
9	5.050	1232	1247	1289	rBV2	189185	469191	100.00%	11.230%
10	5.619	1412	1424	1451	rBV	132896	270029	57.55%	6.463%
11	5.918	1504	1517	1544	rVB5	26038	57140	12.18%	1.368%
12	6.072	1553	1565	1597	rBV	111393	232693	49.59%	5.569%
13	6.989	1840	1850	1877	rBV	41535	79736	16.99%	1.908%
14	7.317	1941	1952	1971	rBV	194311	342571	73.01%	8.199%
15	7.628	2039	2049	2072	rBV	23304	43870	9.35%	1.050%
16	7.937	2136	2145	2149	rBV3	3948	6001	1.28%	0.144%
17	7.979	2149	2158	2172	rVB3	16885	29871	6.37%	0.715%
18	8.088	2182	2192	2235	rBV	215074	425716	90.73%	10.189%
19	8.854	2419	2430	2455	rBV	193174	320258	68.26%	7.665%
20	9.934	2753	2766	2780	rBV2	11275	17485	3.73%	0.418%
21	10.217	2842	2854	2877	rBV	80195	127418	27.16%	3.050%
22	10.358	2886	2898	2916	rBV2	14653	33886	7.22%	0.811%
23	10.741	3007	3017	3027	rBV2	6486	10177	2.17%	0.244%
24	11.088	3119	3125	3136	rVB3	8546	12071	2.57%	0.289%
25	11.249	3165	3175	3200	rBV	179946	284357	60.61%	6.806%
26	11.529	3252	3262	3280	rBV4	20433	41556	8.86%	0.995%
27	11.622	3280	3291	3312	rBV	196258	326942	69.68%	7.825%
28	12.117	3436	3445	3458	rBV4	9764	22877	4.88%	0.548%
29	12.886	3676	3684	3694	rVB3	6099	9002	1.92%	0.215%
30	13.046	3726	3734	3742	rVB3	4667	6453	1.38%	0.154%
31	13.506	3867	3877	3890	rBV2	7511	11572	2.47%	0.277%
32	14.040	4033	4043	4054	rVB2	9587	15252	3.25%	0.365%
33	15.638	4530	4540	4550	rBV2	3433	4845	1.03%	0.116%

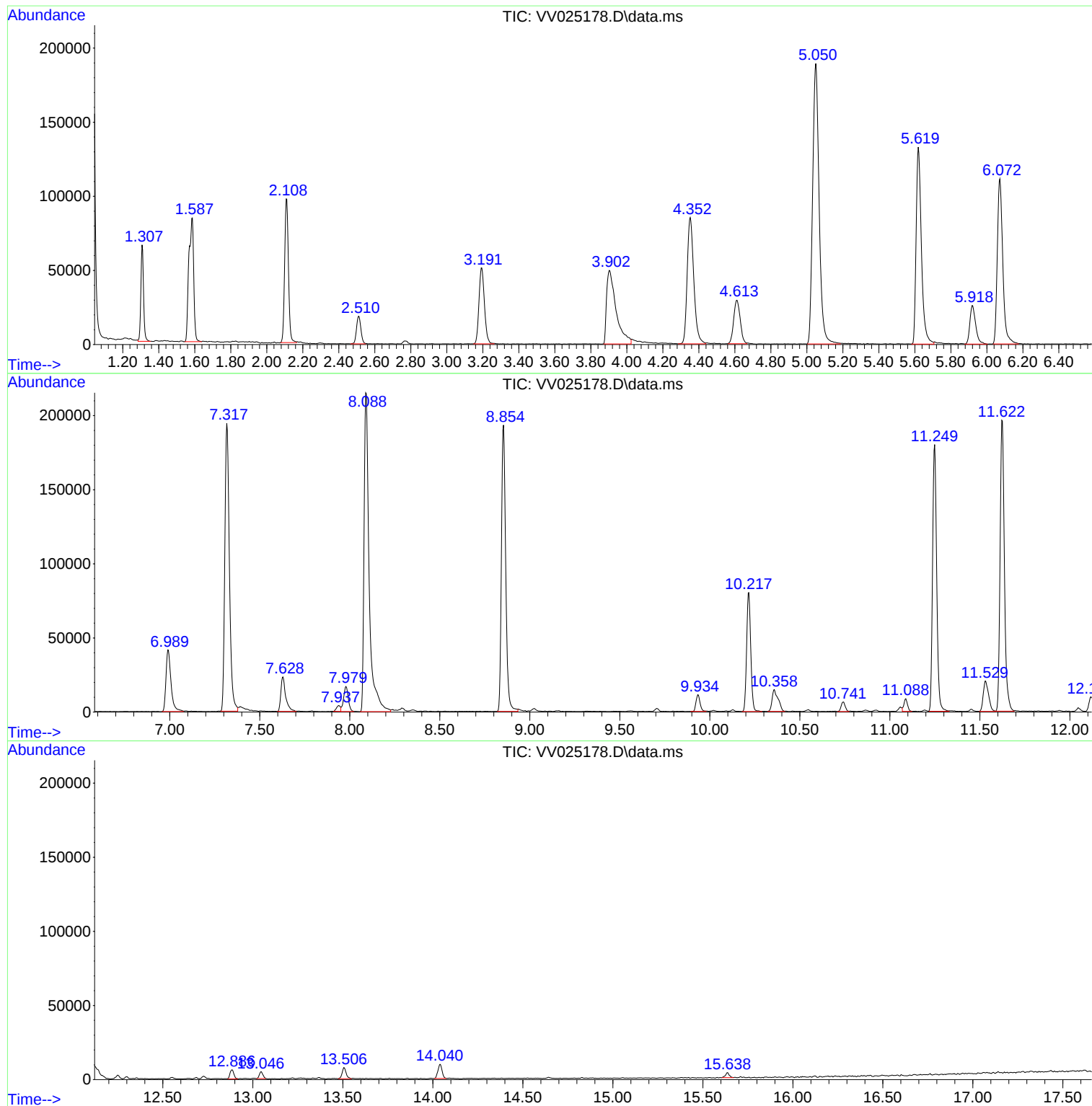
Sum of corrected areas: 4178080

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ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EW607DL

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR031522WMA.M
Quant Title : TRACE VOA SFAM1.0

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



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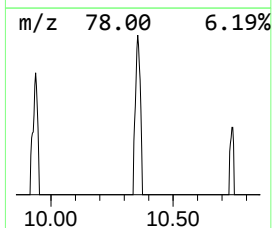
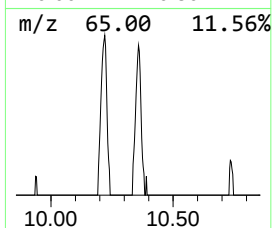
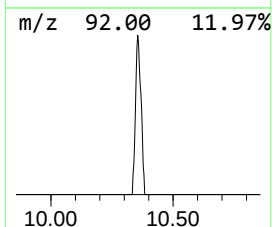
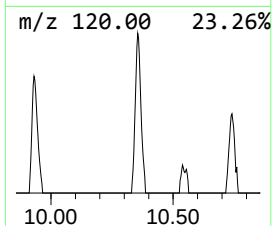
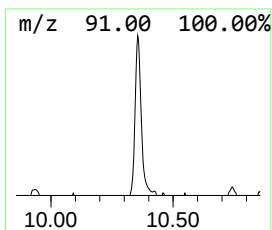
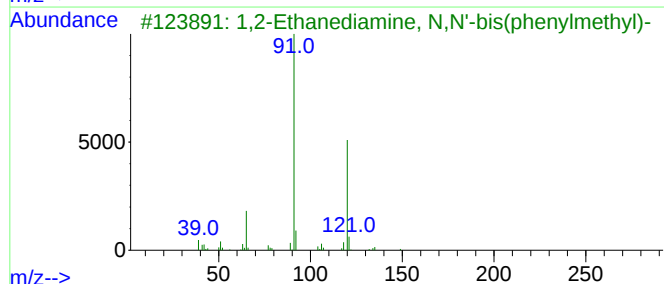
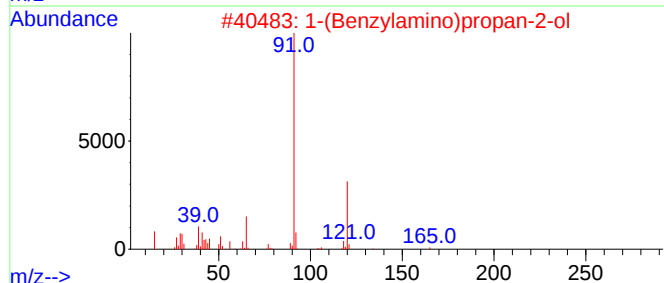
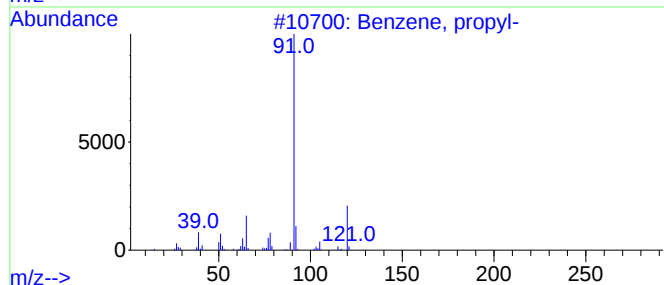
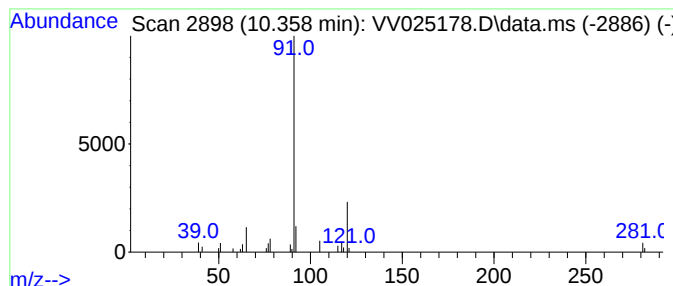
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR031522WMA.M
Quant Title : TRACE VOA SFAM1.0

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

Peak Number 2 Benzene, propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.358	0.60 ug/L	33886	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	81
2			1-(Benzylamino)propan-2-ol	165	C10H15NO	1000486-84-7	72
3			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	72
4			Benzene, propyl-	120	C9H12	000103-65-1	68
5			Benzene, propyl-	120	C9H12	000103-65-1	68



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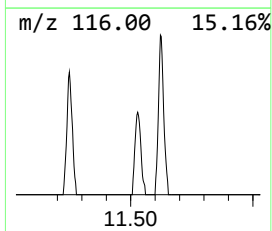
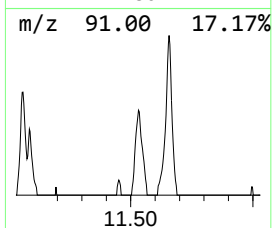
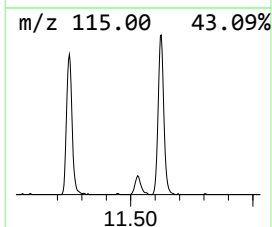
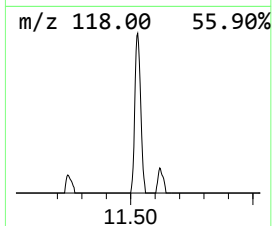
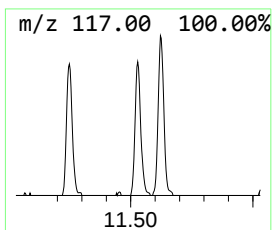
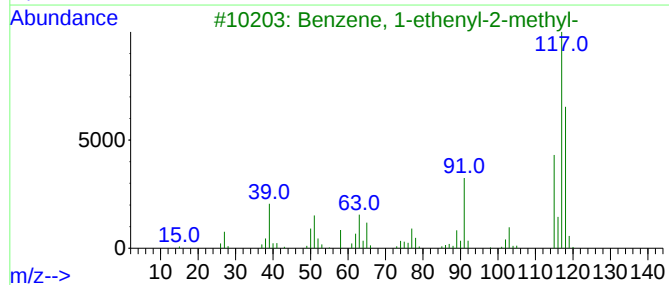
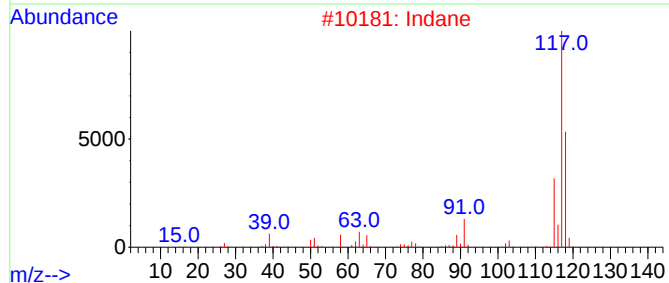
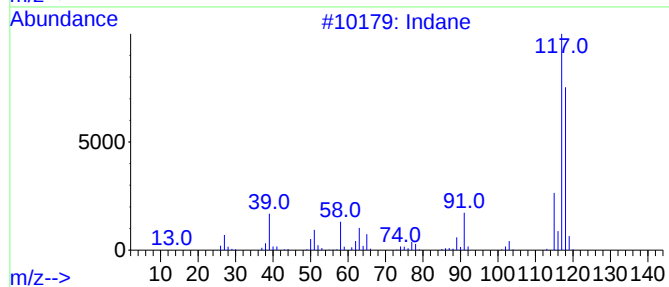
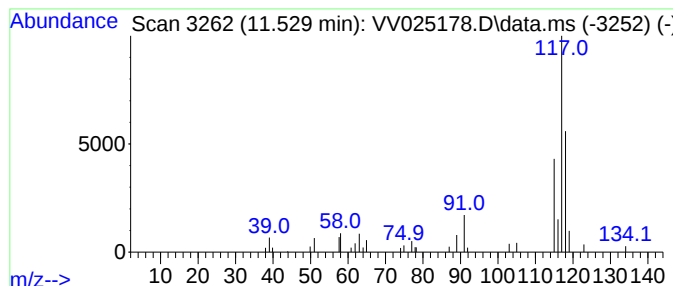
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Quant Title : TRACE VOA SFAM1.0

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

Peak Number 3 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.529	0.73 ug/L	41556	1,4-Dichlorobenzene-d4	11.249

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	76
2	Indane			118	C9H10	000496-11-7	76
3	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	70
4	Benzene, 1-ethenyl-4-methyl-			118	C9H10	000622-97-9	68
5	Indane			118	C9H10	000496-11-7	68



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, propyl-	10.358	0.6	ug/L	33886	3	11.249	284357	5.0
Indane	11.529	0.7	ug/L	41556	3	11.249	284357	5.0