

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091223\
 Data File : BP017141.D
 Acq On : 13 Sep 2023 10:31
 Operator : MA/JU
 Sample : 04324-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHG3

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017141.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.270	25	31	36	rBV	490831	758304	3.15%	0.887%
2	3.317	36	39	43	rVB	65526	85150	0.35%	0.100%
3	3.752	109	113	125	rBV	300722	455135	1.89%	0.532%
4	3.946	141	146	151	rBV	53738	75948	0.32%	0.089%
5	4.587	248	255	282	rVB	14462568	24104553	100.00%	28.202%
6	6.887	643	646	659	rVB9	10363	30691	0.13%	0.036%
7	7.099	676	682	696	rBV	274579	453085	1.88%	0.530%
8	7.293	708	715	723	rBV	1003751	1529560	6.35%	1.790%
9	7.475	736	746	756	rBV	951545	1531578	6.35%	1.792%
10	7.952	820	827	841	rBV	812360	1279285	5.31%	1.497%
11	8.663	940	948	961	rBV	772392	1281996	5.32%	1.500%
12	9.152	1022	1031	1044	rBV	1143302	1838156	7.63%	2.151%
13	9.863	1145	1152	1162	rBV	912882	1447486	6.01%	1.694%
14	10.387	1234	1241	1256	rBV	1320593	2224535	9.23%	2.603%
15	10.775	1299	1307	1316	rBV	1088215	1850049	7.68%	2.165%
16	10.940	1327	1335	1346	rBV	706733	1172944	4.87%	1.372%
17	11.775	1470	1477	1487	rBV	54742	122537	0.51%	0.143%
18	12.357	1568	1576	1582	rVV3	22742	53026	0.22%	0.062%
19	13.457	1756	1763	1778	rBV	516886	845949	3.51%	0.990%
20	14.028	1853	1860	1872	rBV	2732721	3638063	15.09%	4.256%
21	14.310	1901	1908	1926	rBV2	3010114	4301794	17.85%	5.033%
22	14.610	1952	1959	1966	rBV	1748113	2475117	10.27%	2.896%
23	14.834	1991	1997	2010	rBV	156159	313105	1.30%	0.366%
24	15.610	2122	2129	2137	rBV	3905706	5451032	22.61%	6.378%
25	15.734	2144	2150	2165	rVV	1181361	1564906	6.49%	1.831%
26	17.381	2424	2430	2440	rBV	2125857	2925207	12.14%	3.422%
27	17.481	2441	2447	2461	rVV	4412426	5922216	24.57%	6.929%
28	18.004	2532	2536	2541	rBV2	23649	36141	0.15%	0.042%
29	18.222	2569	2573	2580	rBV	94439	130600	0.54%	0.153%
30	19.110	2721	2724	2726	rBV3	28071	32626	0.14%	0.038%
31	19.139	2726	2729	2736	rVB	171347	199171	0.83%	0.233%
32	19.292	2752	2755	2760	rVB2	47754	59520	0.25%	0.070%
33	19.357	2763	2766	2771	rVB	54010	70202	0.29%	0.082%
34	19.457	2779	2783	2788	rBV4	70370	93859	0.39%	0.110%
35	19.722	2822	2828	2839	rBV	5405814	6790051	28.17%	7.944%
36	21.480	3122	3127	3135	rBV	2056141	2666323	11.06%	3.120%

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Integration Parameters: LSCINT.P
Integrator: RTE
Smoothing : OFF Filtering: 5
Sampling : 1 Min Area: 0.1 % of largest Peak
Start Thrs: 0.2 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:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Title : SVOA CALIBRATION

37	23.839	3521	3528	3542	rVB2	2579888	5279480	21.90%	6.177%
38	24.004	3550	3556	3569	rVB	1161110	2382818	9.89%	2.788%

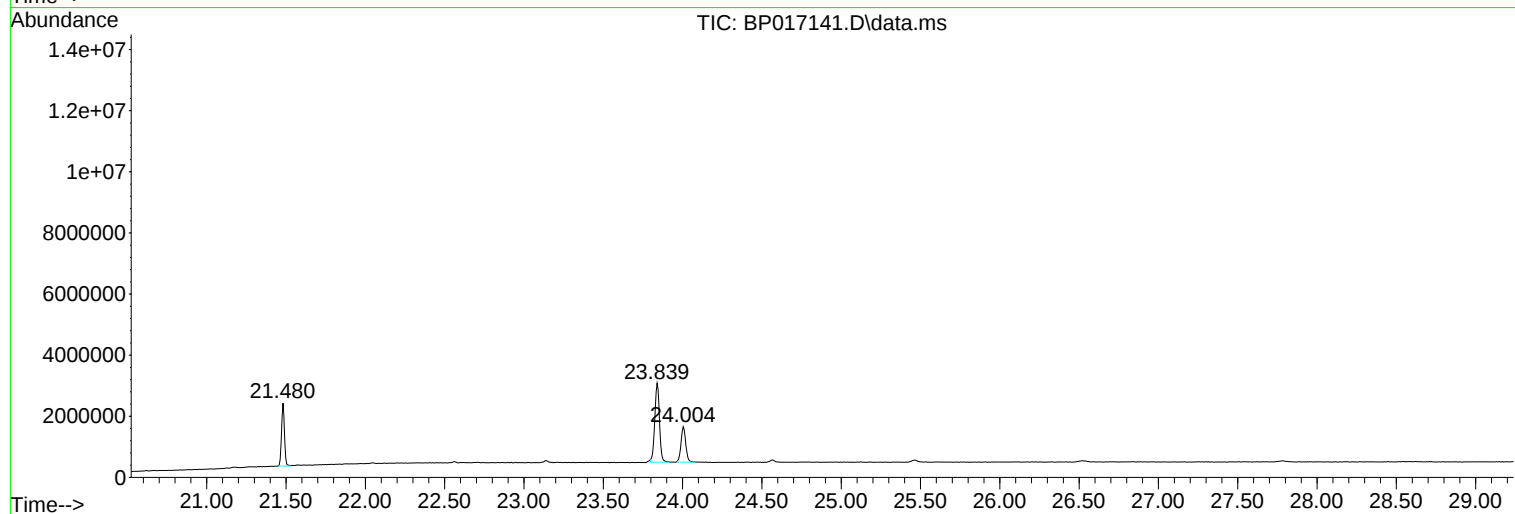
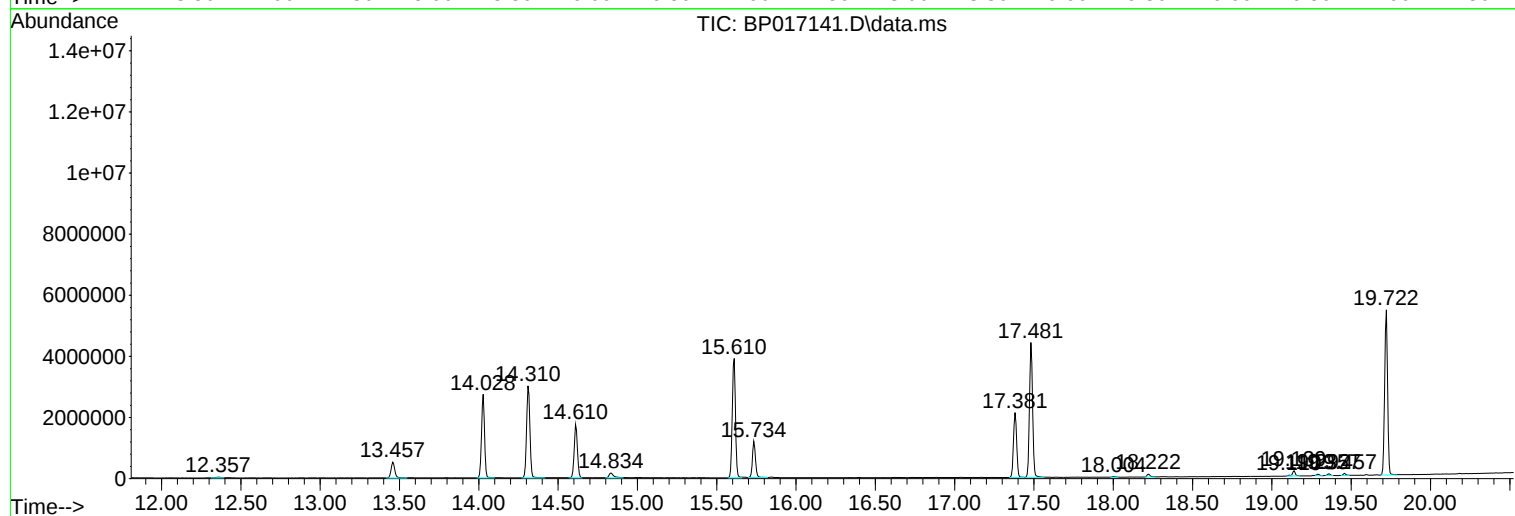
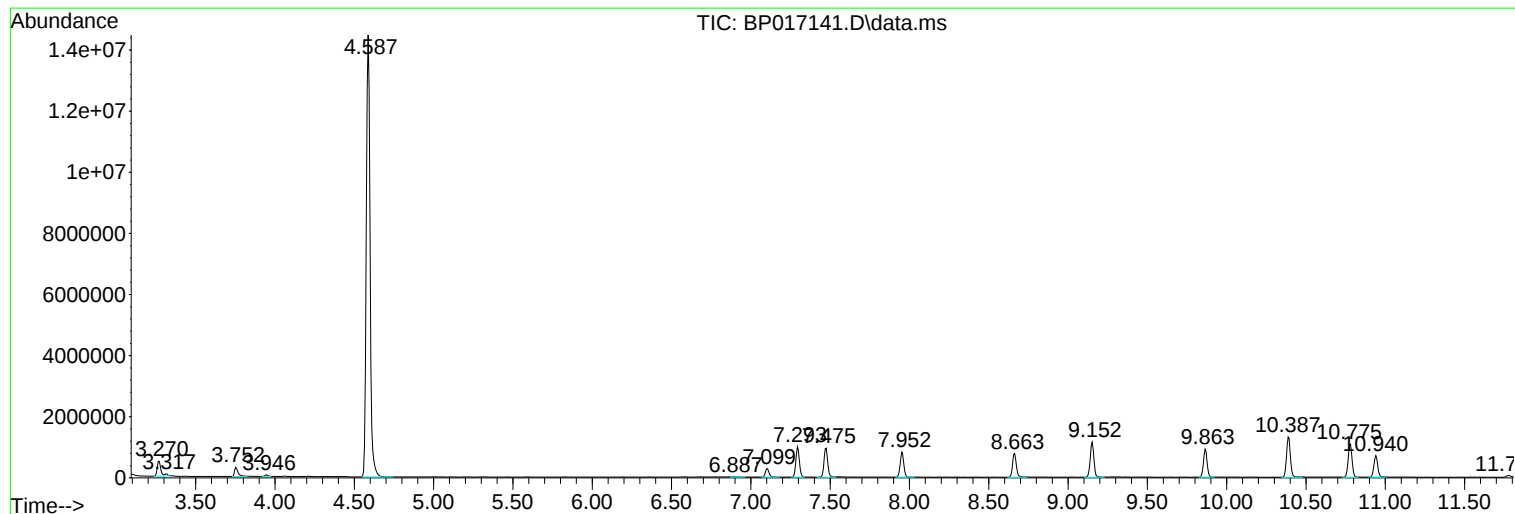
Sum of corrected areas: 85472198

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
Quant Title : SVOA CALIBRATION

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



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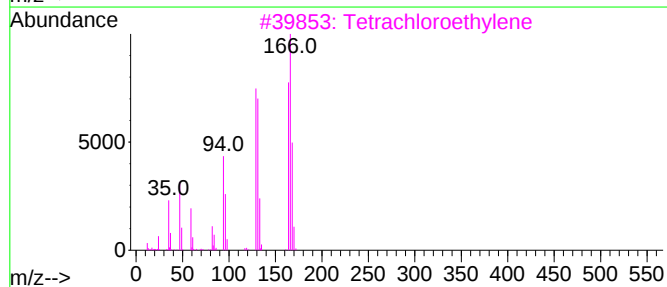
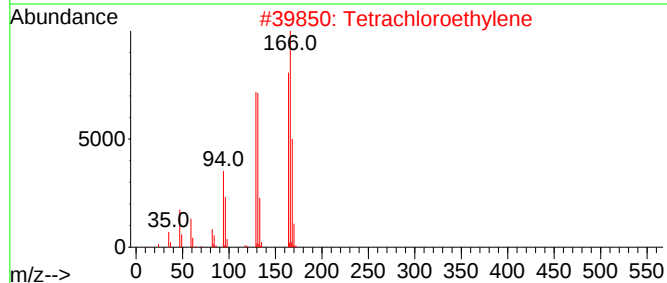
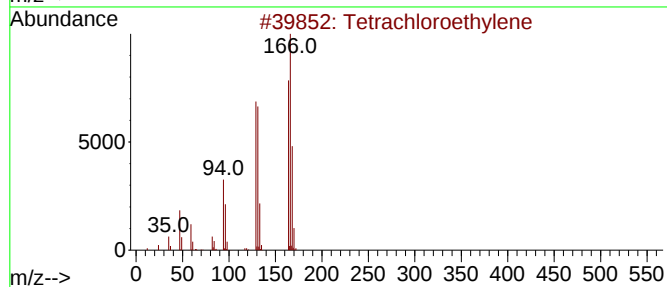
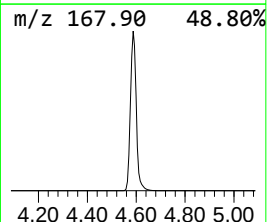
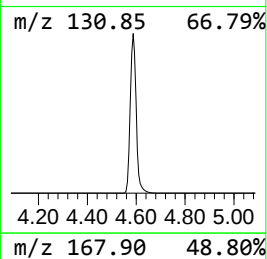
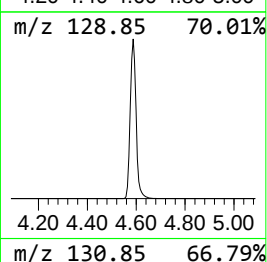
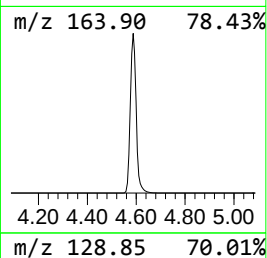
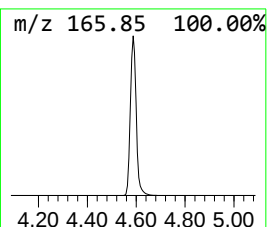
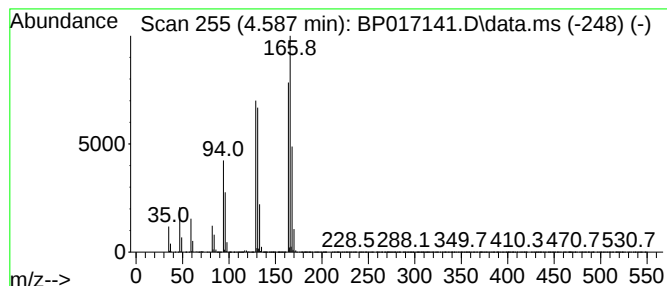
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP091223.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.587	376.84 ng/ul	24104600	1,4-Dichlorobenzene-d4	7.952

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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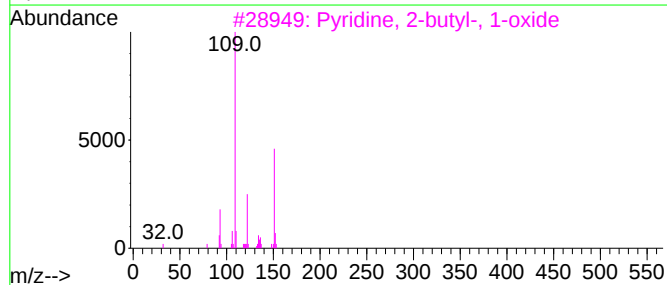
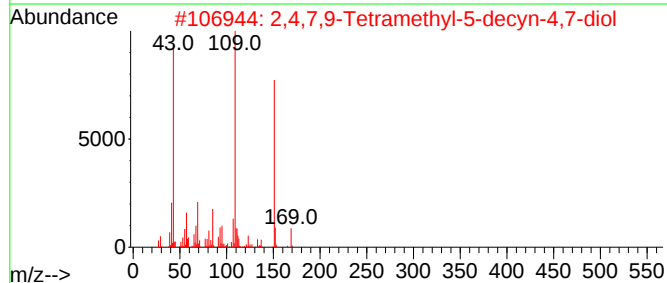
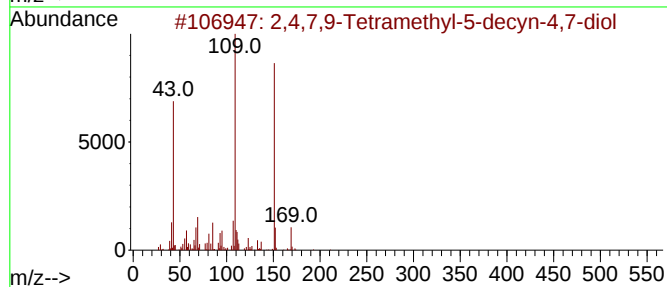
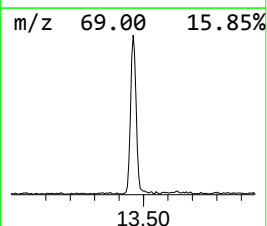
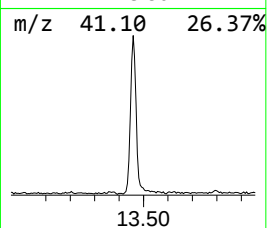
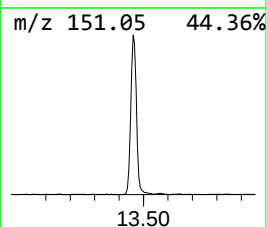
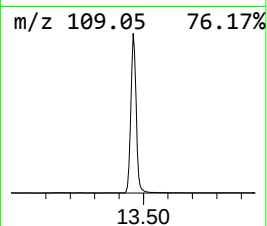
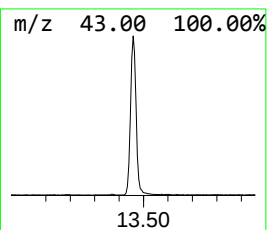
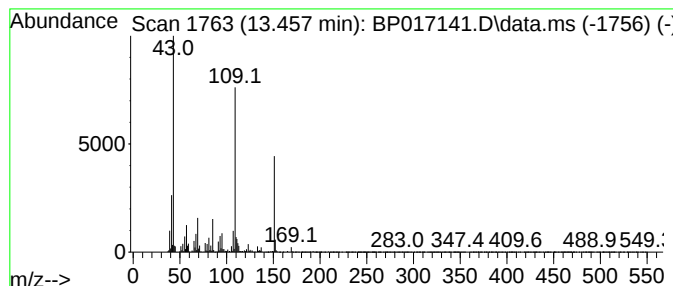
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Peak Number 2 2,4,7,9-Tetramethyl-5-decyn... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.457	6.84 ng/ul	845949	Acenaphthene-d10	14.610

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	90		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	64		
3	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	59		
4	Diacetamate	193	C10H11NO3	002623-33-8	59		
5	3-Pyridinol, 4-methyl-, acetate ...	151	C8H9NO2	001006-96-8	59		



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.587	376.8	ng/ul	24104600	1	7.952	1279290	20.0
2,4,7,9-Tetrame...	13.457	6.8	ng/ul	845949	3	14.610	2475120	20.0