

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111324\
 Data File : BP023137.D
 Acq On : 15 Nov 2024 13:23
 Operator : RC/JU
 Sample : P4740-02
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CC3

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-BP110924.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP023137.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.134	22	25	31	rVB	12725	15252	0.11%	0.046%
2	3.564	95	98	106	rBV	27734	51182	0.38%	0.155%
3	4.369	228	235	255	rBV	9458084	13443749	100.00%	40.693%
4	4.934	326	331	337	rBV3	11526	20565	0.15%	0.062%
5	6.793	641	647	664	rBV2	40548	104134	0.77%	0.315%
6	6.928	664	670	683	rVB	156942	263418	1.96%	0.797%
7	7.128	697	704	720	rBV	193291	338696	2.52%	1.025%
8	7.575	774	780	791	rBV	221059	378599	2.82%	1.146%
9	8.293	896	902	923	rBV2	121142	250014	1.86%	0.757%
10	8.722	967	975	989	rBV	205016	393752	2.93%	1.192%
11	9.434	1087	1096	1118	rBV2	172444	359254	2.67%	1.087%
12	9.981	1181	1189	1217	rBV	228588	547852	4.08%	1.658%
13	10.328	1240	1248	1258	rBV	319161	523808	3.90%	1.586%
14	10.481	1267	1274	1292	rBV	133046	300344	2.23%	0.909%
15	11.181	1384	1393	1408	rBV5	11151	30854	0.23%	0.093%
16	13.087	1711	1717	1727	rBV2	53138	99266	0.74%	0.300%
17	13.598	1794	1804	1823	rBV	788957	1001502	7.45%	3.031%
18	13.887	1844	1853	1870	rBV2	620402	995952	7.41%	3.015%
19	14.198	1897	1906	1916	rBV	512969	798061	5.94%	2.416%
20	14.510	1953	1959	1964	rBV3	22531	57202	0.43%	0.173%
21	15.210	2069	2078	2091	rBV	769422	1466726	10.91%	4.440%
22	15.363	2098	2104	2116	rBV	292589	516737	3.84%	1.564%
23	15.451	2116	2119	2125	rVB7	8631	14025	0.10%	0.042%
24	16.928	2364	2370	2375	rBV2	35589	46778	0.35%	0.142%
25	17.004	2375	2383	2392	rBV	576274	1068758	7.95%	3.235%
26	17.110	2394	2401	2415	rVV	1204152	1835889	13.66%	5.557%
27	17.945	2536	2543	2551	rBV2	55279	90281	0.67%	0.273%
28	18.380	2614	2617	2622	rVB5	17982	22854	0.17%	0.069%
29	18.933	2708	2711	2719	rVB9	13069	16079	0.12%	0.049%
30	19.039	2725	2729	2733	rBV2	19703	31979	0.24%	0.097%
31	19.116	2738	2742	2746	rVB	11954	19301	0.14%	0.058%
32	19.280	2766	2770	2777	rBV6	18452	27665	0.21%	0.084%
33	19.492	2800	2806	2816	rBV	1792505	2391942	17.79%	7.240%
34	21.445	3132	3138	3148	rVB	753119	1430076	10.64%	4.329%
35	24.445	3638	3648	3662	rVB	952058	2559386	19.04%	7.747%
36	24.662	3677	3685	3703	rVB	516708	1525262	11.35%	4.617%

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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-BP110924.MA.M
Title : SVOA CALIBRATION

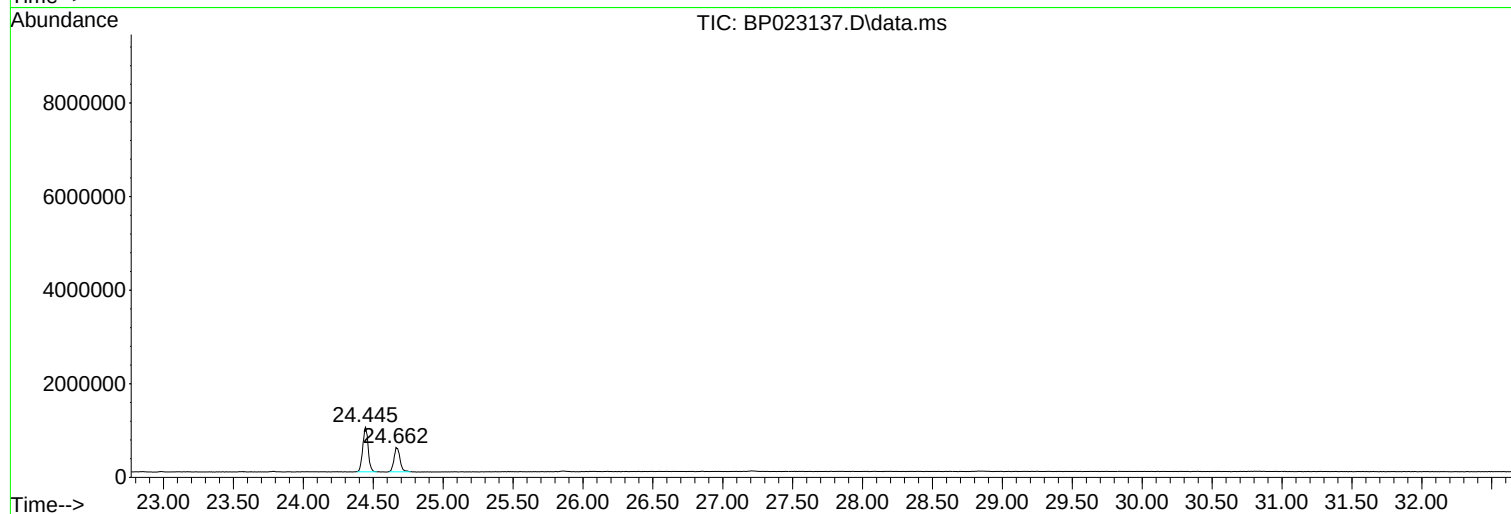
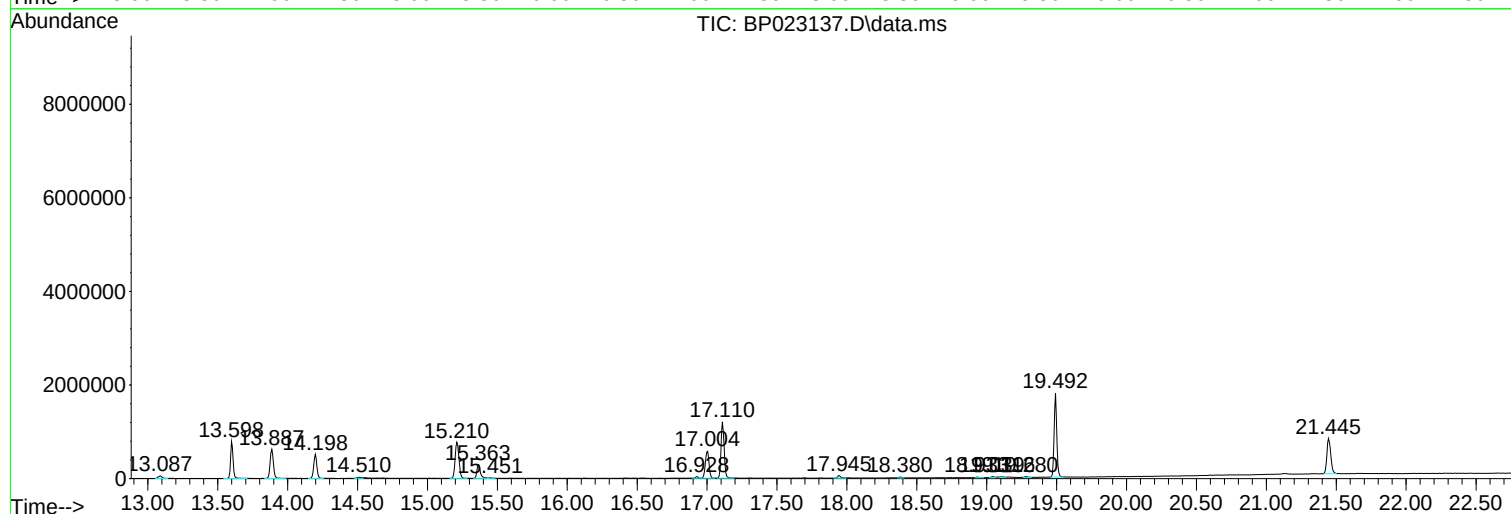
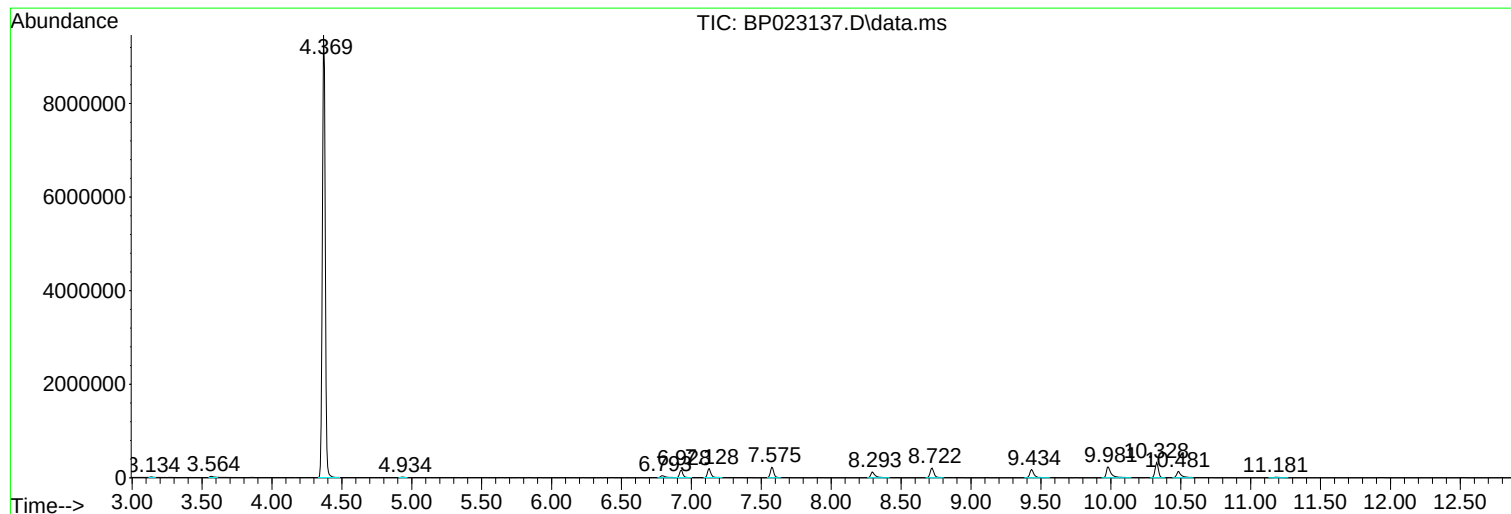
Sum of corrected areas: 33037194

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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



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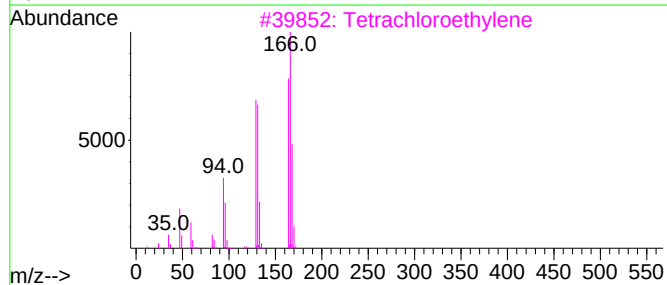
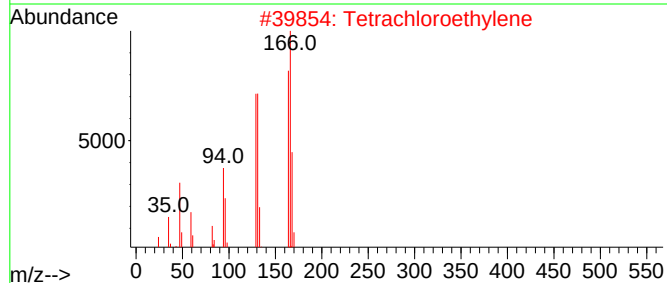
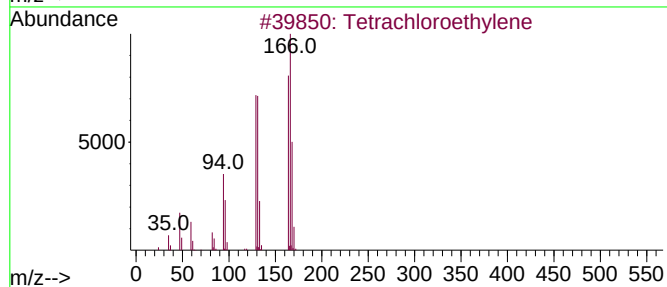
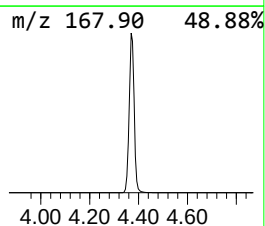
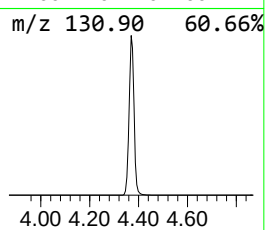
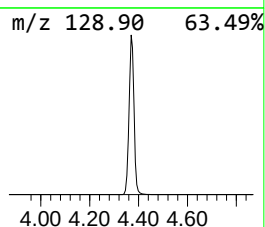
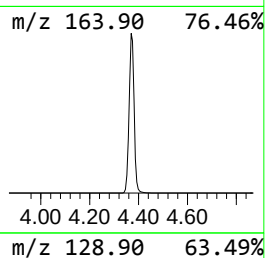
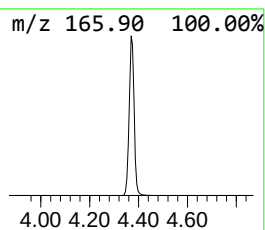
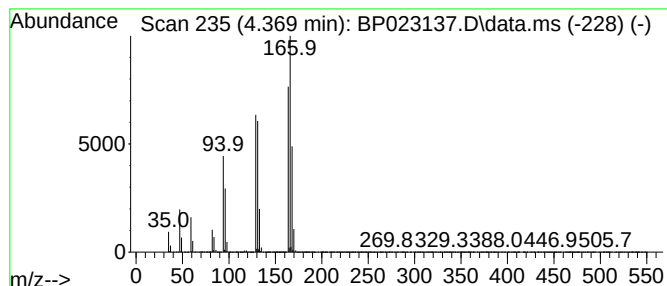
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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.
4.369	710.18 ng/ul	13443700	1,4-Dichlorobenzene-d4	7.575

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	99
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	94
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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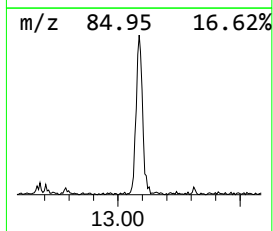
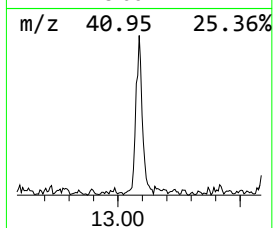
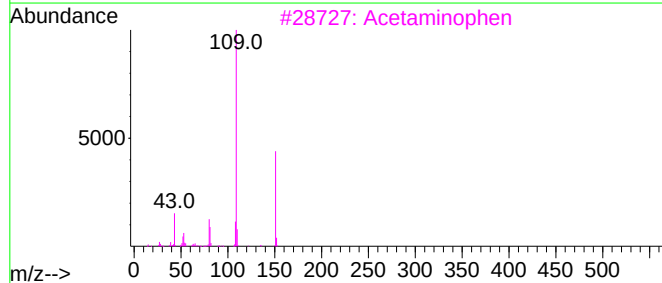
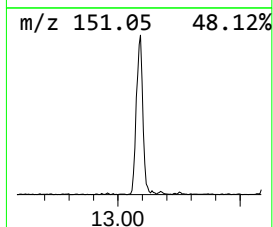
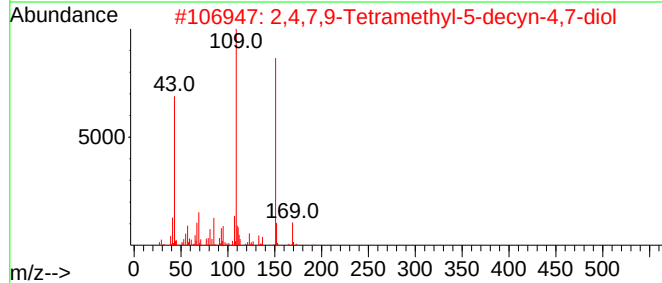
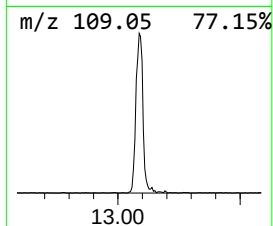
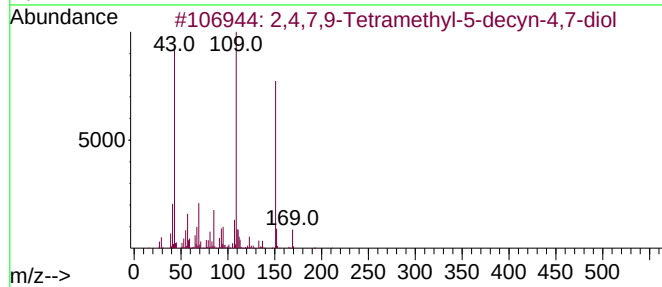
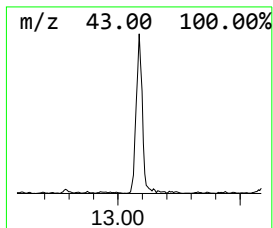
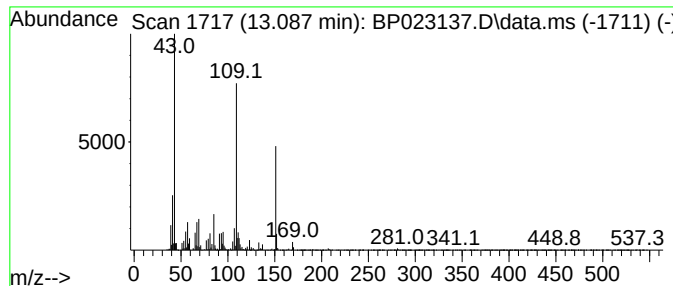
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION

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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.087	2.49 ng/ul	99266	Acenaphthene-d10	14.198

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
2	2,4,7,9-Tetramethyl-5-decyn-4,7-...	226	C14H26O2	000126-86-3	91		
3	Acetaminophen	151	C8H9NO2	000103-90-2	59		
4	Metacetamol	151	C8H9NO2	000621-42-1	59		
5	Acetaminophen	151	C8H9NO2	000103-90-2	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...	4.369	710.2	ng/ul	13443700	1	7.575	378599	20.0
2,4,7,9-Tetrame...	13.087	2.5	ng/ul	99266	3	14.198	798061	20.0