

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017504.D
 Acq On : 03 Oct 2023 07:01
 Operator : MA/JU
 Sample : 04588-06
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJE3

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-BP092023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017504.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.240	20	26	45	rVB	2732835	4113713	9.58%	4.659%
2	3.722	105	108	122	rBV	238515	372819	0.87%	0.422%
3	4.552	242	249	253	rBV	21740731	42952594	100.00%	48.646%
4	5.011	323	327	333	rVB	33366	49708	0.12%	0.056%
5	6.363	552	557	562	rBV	30819	43390	0.10%	0.049%
6	7.093	674	681	693	rBV	154002	270641	0.63%	0.307%
7	7.275	705	712	726	rVB	693494	1063492	2.48%	1.204%
8	7.457	734	743	755	rBV	654576	1080196	2.51%	1.223%
9	7.934	816	824	833	rBV	556027	855531	1.99%	0.969%
10	8.657	940	947	958	rBV	475887	774722	1.80%	0.877%
11	9.146	1019	1030	1044	rBV	800044	1365339	3.18%	1.546%
12	9.857	1144	1151	1164	rBV	633496	1036102	2.41%	1.173%
13	10.393	1229	1242	1262	rBV	861965	1539443	3.58%	1.743%
14	10.775	1299	1307	1316	rBV	721155	1201216	2.80%	1.360%
15	10.946	1328	1336	1349	rBV	761187	1321952	3.08%	1.497%
16	12.363	1568	1577	1585	rVB4	18025	44990	0.10%	0.051%
17	13.469	1758	1765	1778	rBV	372354	586825	1.37%	0.665%
18	14.045	1857	1863	1875	rBV	1710108	2267222	5.28%	2.568%
19	14.328	1904	1911	1922	rBV	1914172	2755772	6.42%	3.121%
20	14.634	1956	1963	1971	rBV	1008509	1486941	3.46%	1.684%
21	14.898	2002	2008	2020	rBV	48498	120690	0.28%	0.137%
22	15.634	2126	2133	2143	rBV	2377235	3318835	7.73%	3.759%
23	15.775	2151	2157	2166	rBV	477813	644667	1.50%	0.730%
24	17.422	2430	2437	2447	rBV	1220282	1685923	3.93%	1.909%
25	17.522	2448	2454	2465	rBV	2665219	3630623	8.45%	4.112%
26	18.269	2577	2581	2594	rBV2	83780	160493	0.37%	0.182%
27	19.186	2734	2737	2746	rVB3	51513	67725	0.16%	0.077%
28	19.416	2772	2776	2779	rVB2	52652	66696	0.16%	0.076%
29	19.516	2789	2793	2805	rBV3	66110	114404	0.27%	0.130%
30	19.774	2832	2837	2846	rBV	3542163	4548576	10.59%	5.151%
31	21.557	3135	3140	3147	rBV	1459979	1896180	4.41%	2.148%
32	23.974	3543	3551	3566	rVB	2296494	4837037	11.26%	5.478%
33	24.145	3573	3580	3591	rVB	944466	2021711	4.71%	2.290%

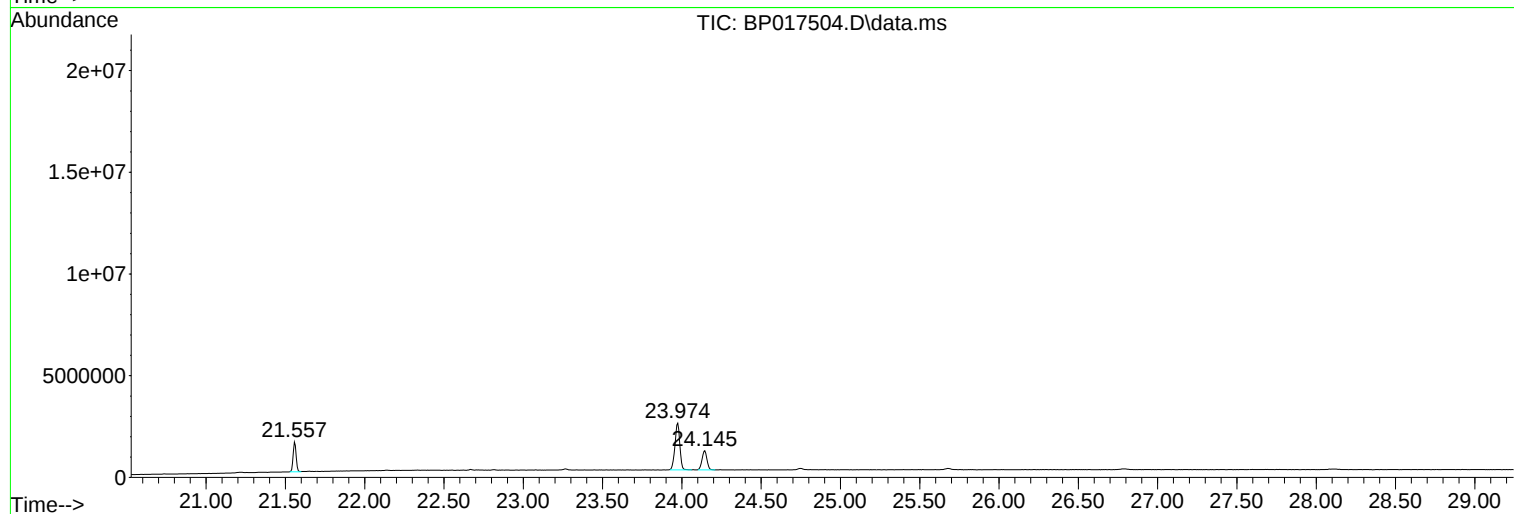
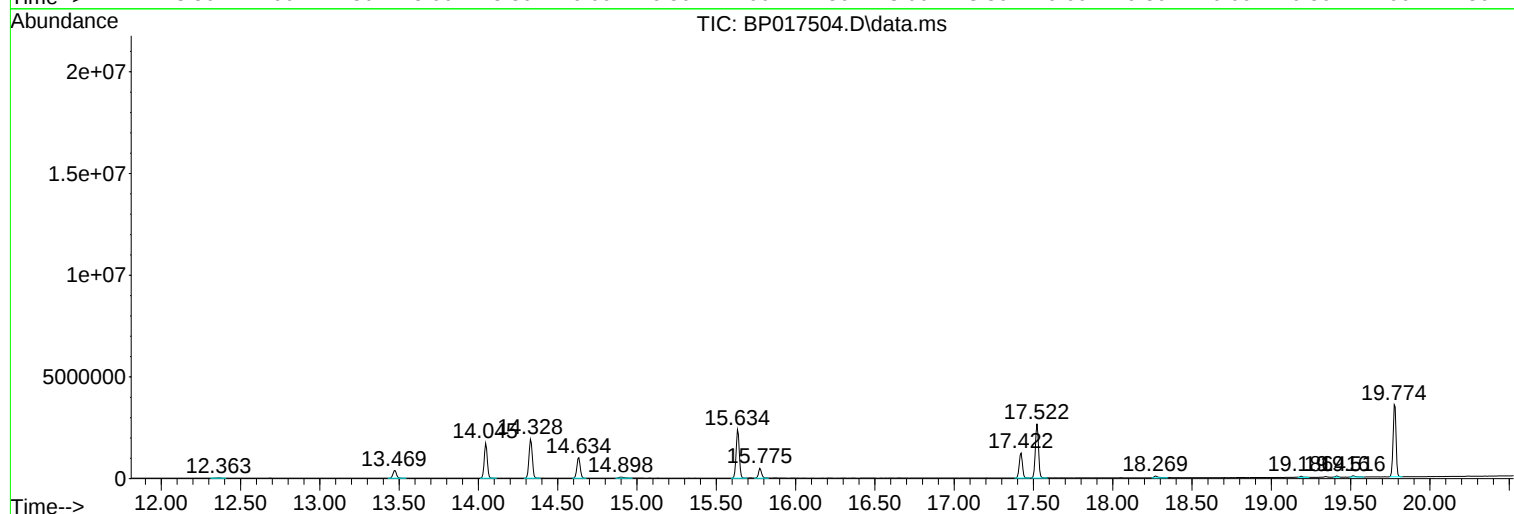
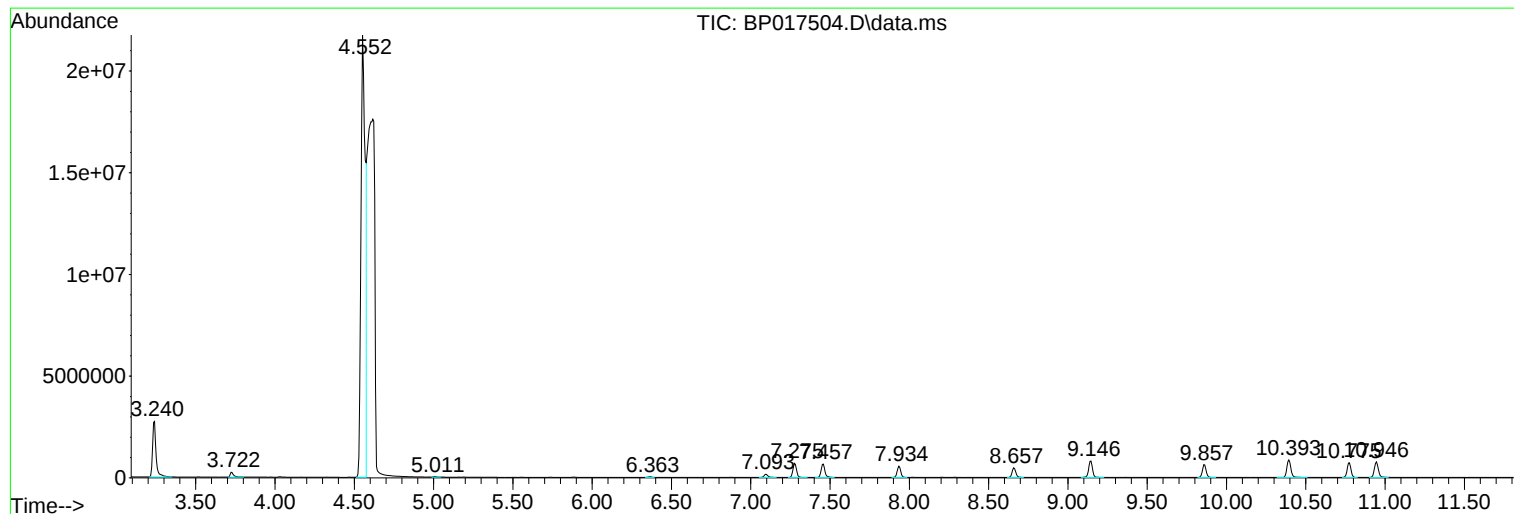
Sum of corrected areas: 88296168

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



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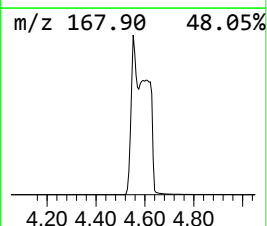
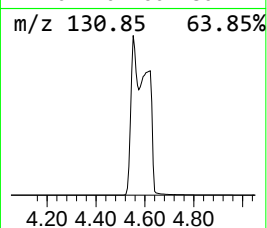
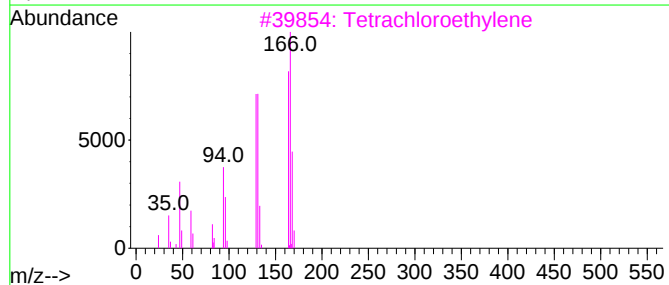
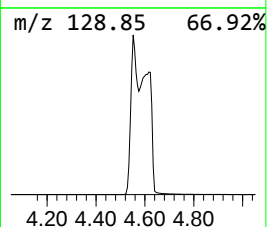
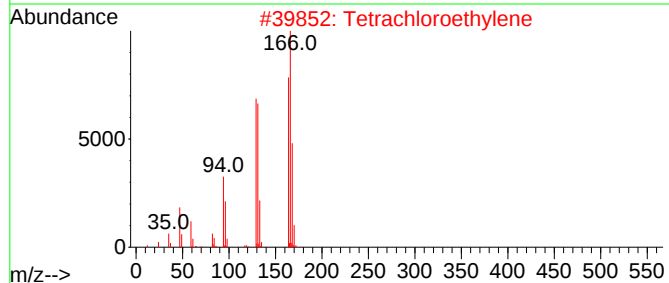
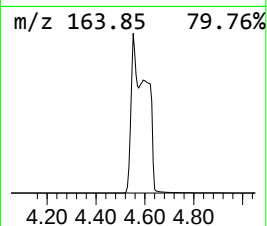
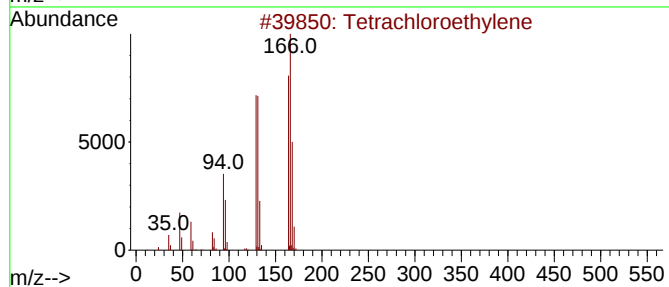
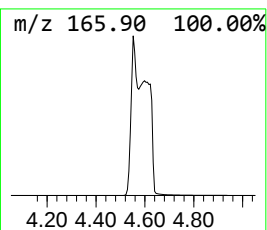
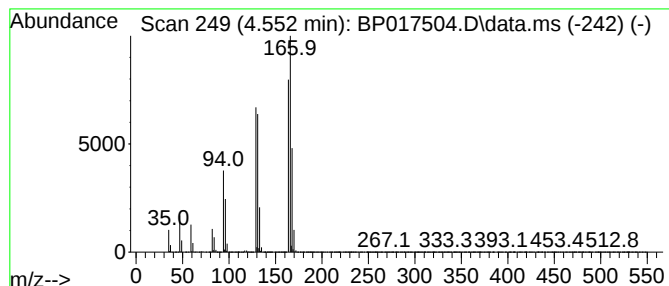
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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.
4.552	1004.12 ng/ul	42952600	1,4-Dichlorobenzene-d4	7.934

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	97
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
5			2,4-Dichloro-6-fluoropyrimidine	166	C4HC12FN2	003833-57-6	38



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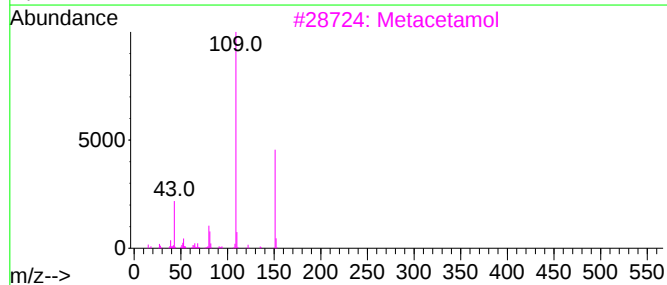
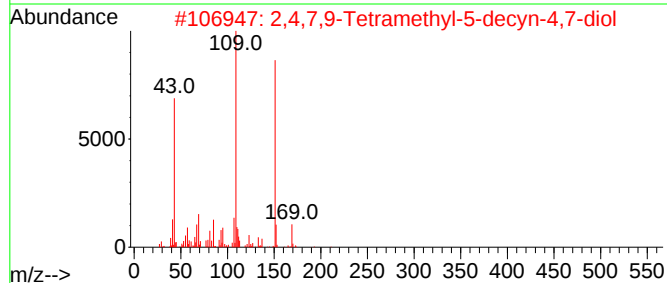
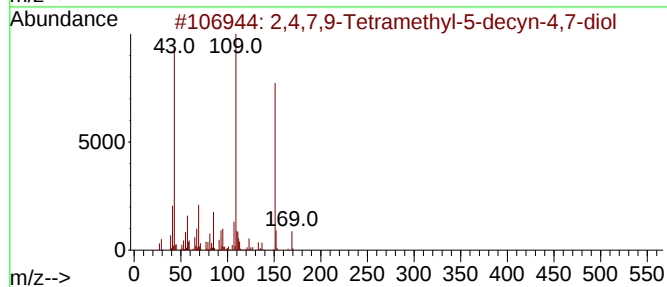
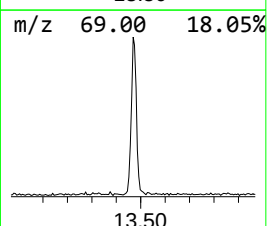
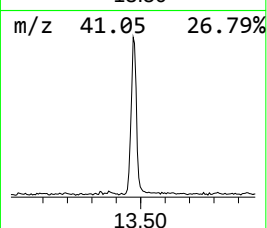
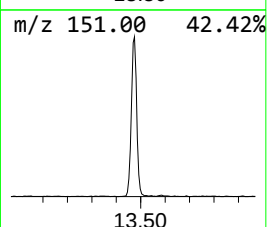
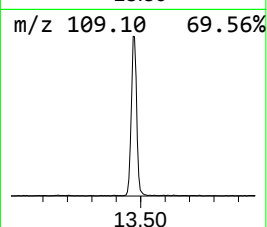
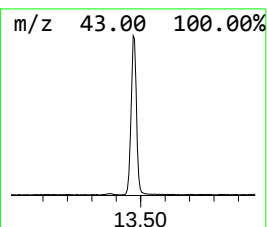
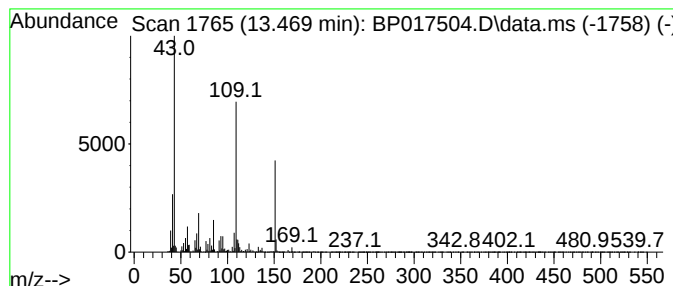
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.469	7.89 ng/ul	586825	Acenaphthene-d10	14.634	
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	83
3	Metacetamol	151	C8H9NO2	000621-42-1	58
4	p-Isopropoxyaniline	151	C9H13NO	007664-66-6	53
5	Pyridine, 2-butyl-, 1-oxide	151	C9H13NO	031396-32-4	53



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Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Tetrachloroethy...	4.552	1004.1	ng/ul	42952600	1	7.934	855531	20.0
2,4,7,9-Tetrame...	13.469	7.9	ng/ul	586825	3	14.634	1486940	20.0