

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017505.D
 Acq On : 03 Oct 2023 07:35
 Operator : MA/JU
 Sample : 04588-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBJM4

Integration Parameters: LSCINT.P

Integrator: RTE

Soothing : 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: BP017505.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.234	20	25	47	rVB	2869135	4205028	9.80%	4.688%
2	3.728	105	109	124	rBV	181550	290990	0.68%	0.324%
3	4.552	242	249	253	rBV	21743391	42917600	100.00%	47.850%
4	5.011	323	327	337	rVB	37378	62066	0.14%	0.069%
5	6.363	552	557	565	rVB	34595	50676	0.12%	0.056%
6	7.093	674	681	695	rVB	138028	235180	0.55%	0.262%
7	7.275	703	712	721	rBV	626935	950146	2.21%	1.059%
8	7.452	734	742	755	rBV	573318	957028	2.23%	1.067%
9	7.934	816	824	833	rBV	575848	886484	2.07%	0.988%
10	8.657	940	947	959	rBV	406255	675530	1.57%	0.753%
11	9.140	1018	1029	1041	rBV	731608	1218007	2.84%	1.358%
12	9.857	1144	1151	1164	rBV	573178	934598	2.18%	1.042%
13	10.393	1235	1242	1263	rBV	817352	1393731	3.25%	1.554%
14	10.775	1298	1307	1320	rBV	740951	1252310	2.92%	1.396%
15	10.946	1328	1336	1351	rBV	683571	1178111	2.75%	1.314%
16	13.469	1759	1765	1777	rBV	326724	529951	1.23%	0.591%
17	14.045	1857	1863	1877	rBV	1737209	2307931	5.38%	2.573%
18	14.328	1904	1911	1922	rBV2	1910806	2721588	6.34%	3.034%
19	14.634	1954	1963	1971	rBV	1108494	1617712	3.77%	1.804%
20	14.898	2002	2008	2021	rBV2	56469	139815	0.33%	0.156%
21	15.633	2127	2133	2142	rBV	2524223	3464961	8.07%	3.863%
22	15.775	2152	2157	2170	rBV	526611	721764	1.68%	0.805%
23	17.422	2431	2437	2445	rBV	1419077	1928879	4.49%	2.151%
24	17.522	2448	2454	2472	rVV	2937270	4041880	9.42%	4.506%
25	18.269	2577	2581	2591	rBV2	58742	99022	0.23%	0.110%
26	19.192	2734	2738	2747	rVB2	75744	100367	0.23%	0.112%
27	19.345	2760	2764	2769	rVB3	37599	51131	0.12%	0.057%
28	19.416	2771	2776	2780	rBV2	46941	65641	0.15%	0.073%
29	19.516	2789	2793	2800	rBV9	31159	55357	0.13%	0.062%
30	19.774	2831	2837	2847	rBV	3732104	4996095	11.64%	5.570%
31	21.557	3135	3140	3147	rBV	1744684	2186284	5.09%	2.438%
32	23.974	3542	3551	3562	rVB2	2416787	5119335	11.93%	5.708%
33	24.139	3572	3579	3592	rVB	1112587	2336870	5.45%	2.605%

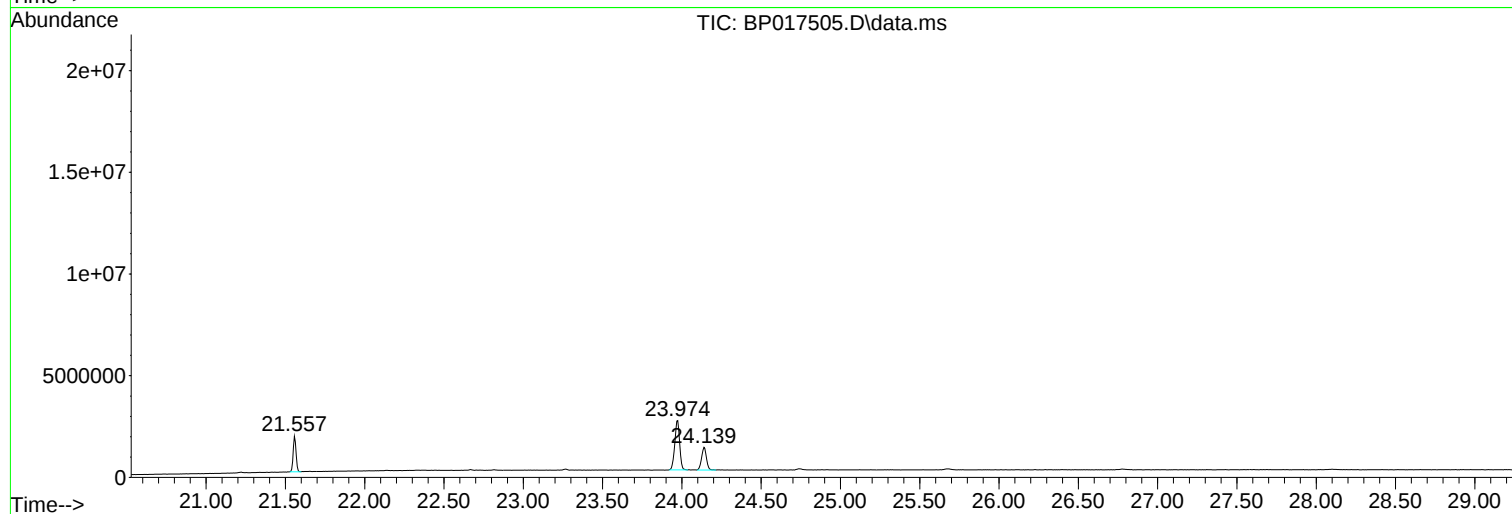
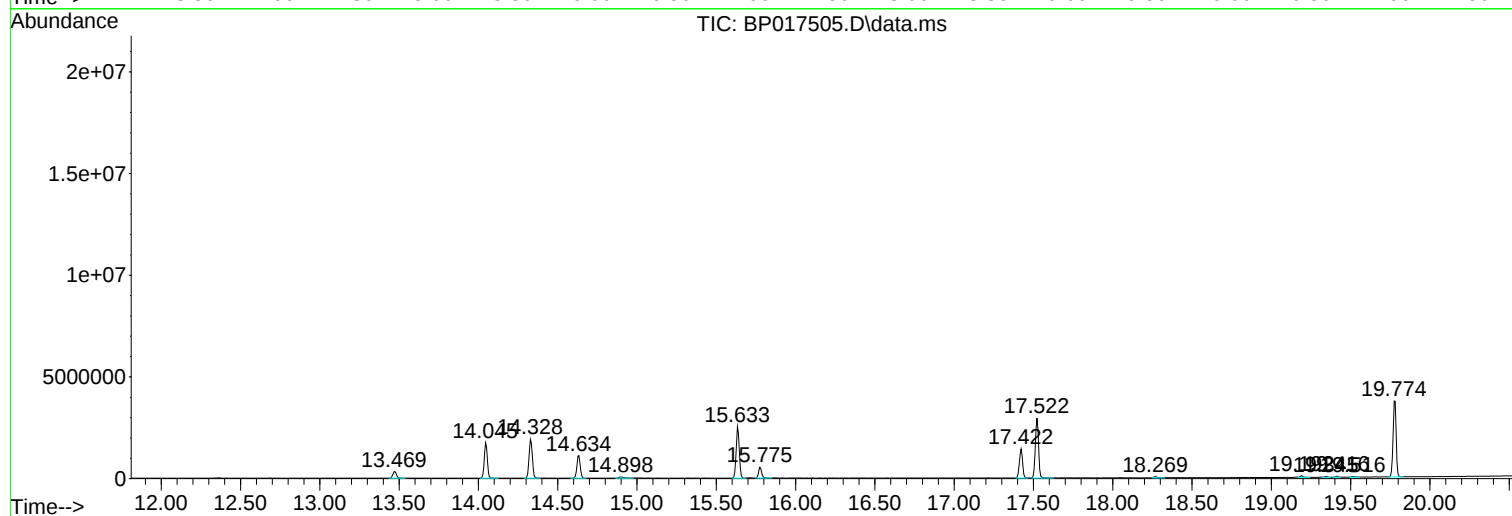
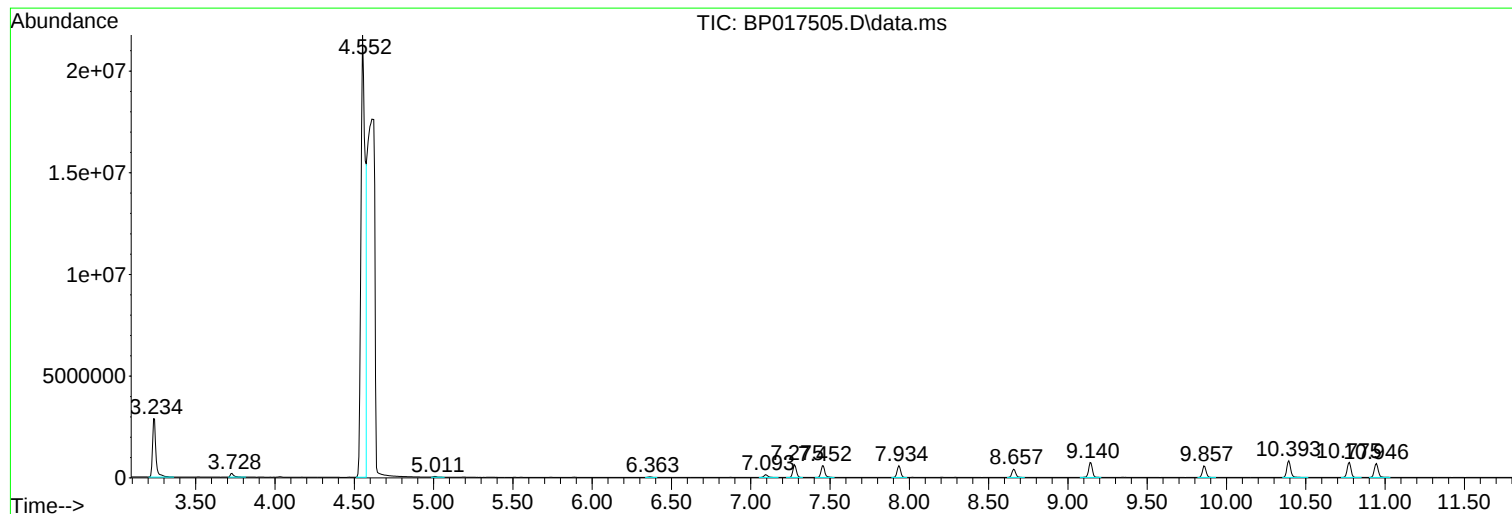
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



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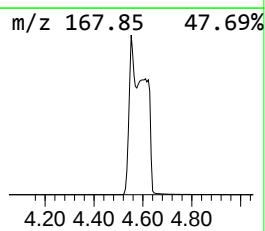
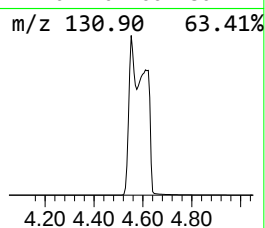
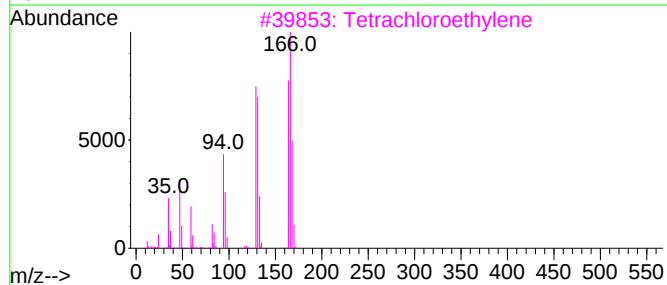
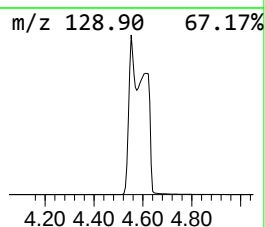
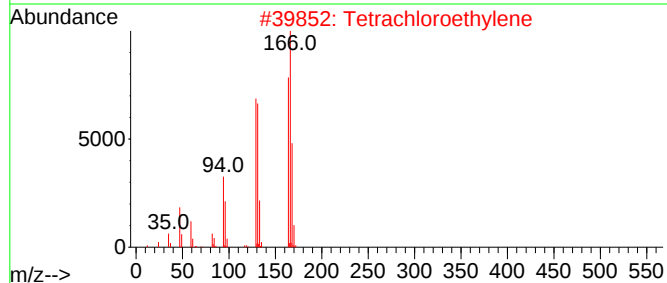
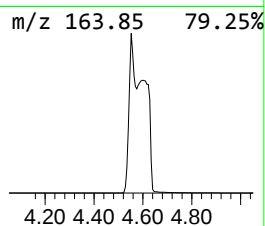
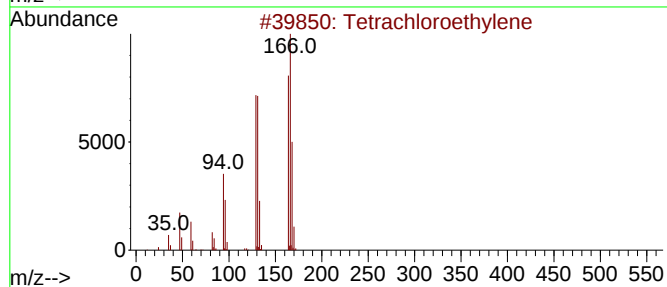
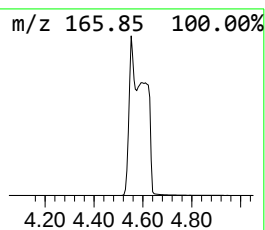
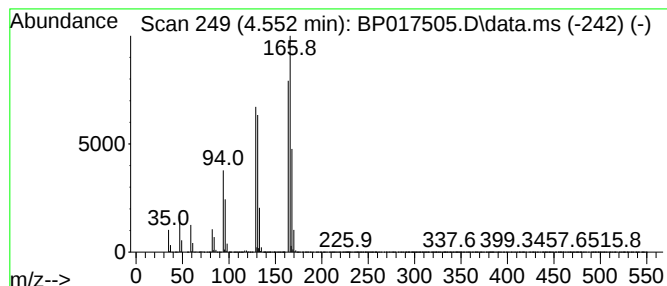
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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	968.27 ng/ul	42917600	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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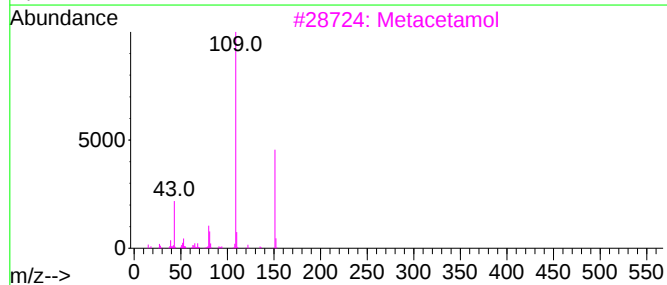
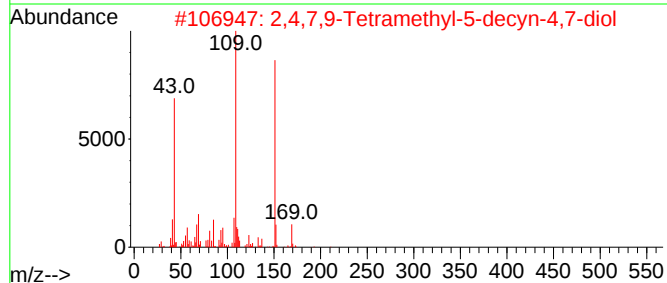
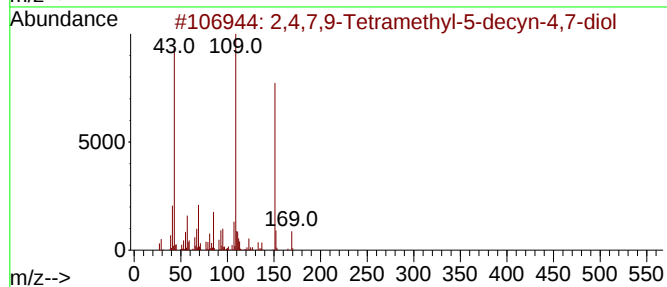
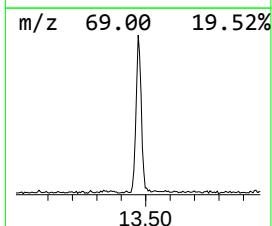
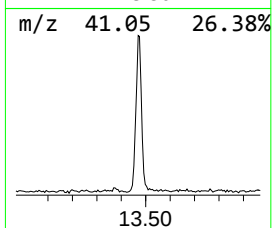
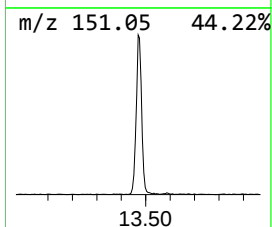
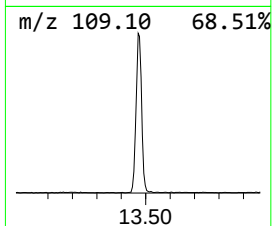
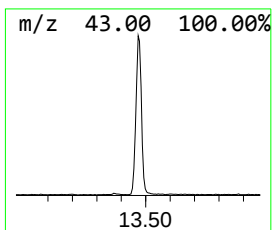
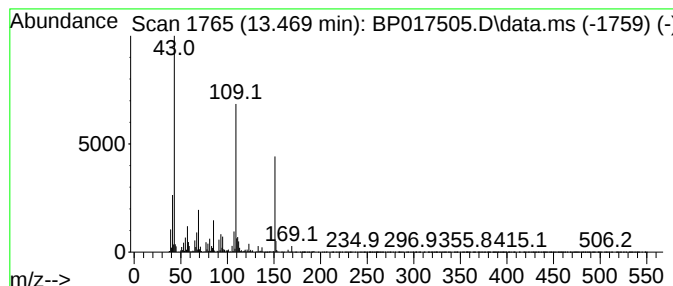
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Quant Title : SVOA CALIBRATION

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	6.55 ng/ul	529951	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	91
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	Butanamide, N-acetyl-N-(4-hydrox...		221	C12H15NO3	1000107-08-7	53
5	3-Pyridinol, 4-methyl-, acetate ...		151	C8H9NO2	001006-96-8	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	968.3	ng/ul	42917600	1	7.934	886484	20.0
2,4,7,9-Tetrame...	13.469	6.5	ng/ul	529951	3	14.634	1617710	20.0