

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082323\
 Data File : BP016716.D
 Acq On : 23 Aug 2023 15:50
 Operator : MA/JU
 Sample : 04105-03
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHAB8

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-BP081423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016716.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.340	5	9	14	rBV	253130	396766	4.08%	0.423%
2	3.387	14	17	21	rVV2	118907	178842	1.84%	0.191%
3	3.423	21	23	31	rVB	66340	89117	0.92%	0.095%
4	3.828	88	92	107	rBV	367938	607939	6.25%	0.649%
5	4.028	121	126	132	rVB2	42594	65417	0.67%	0.070%
6	4.146	141	146	155	rVB	20407	37103	0.38%	0.040%
7	4.528	209	211	217	rVB4	10052	15077	0.15%	0.016%
8	4.670	229	235	248	rVB	1467171	2223473	22.86%	2.373%
9	5.081	302	305	310	rVB4	7674	11913	0.12%	0.013%
10	6.687	572	578	583	rBV7	5054	10007	0.10%	0.011%
11	7.181	656	662	670	rBV	348189	540332	5.55%	0.577%
12	7.375	686	695	708	rBV	1485512	2319482	23.84%	2.476%
13	7.558	719	726	736	rBV	1306366	2044026	21.01%	2.182%
14	7.769	758	762	766	rVB7	8960	14393	0.15%	0.015%
15	8.040	801	808	819	rBV	1181967	1855064	19.07%	1.980%
16	8.740	920	927	938	rBV	1004321	1645808	16.92%	1.757%
17	8.828	938	942	952	rVB	27695	53528	0.55%	0.057%
18	9.240	1004	1012	1022	rBV	1751093	2871108	29.51%	3.064%
19	9.952	1124	1133	1146	rBV	1316297	2177295	22.38%	2.324%
20	10.475	1215	1222	1244	rBV	1776419	3130014	32.17%	3.341%
21	10.869	1282	1289	1297	rBV	1735634	2876282	29.57%	3.070%
22	11.028	1307	1316	1327	rBV	1955456	3263963	33.55%	3.484%
23	12.446	1551	1557	1565	rBV	34819	62451	0.64%	0.067%
24	13.534	1738	1742	1747	rVV5	12819	20065	0.21%	0.021%
25	13.657	1758	1763	1768	rBV7	6223	10825	0.11%	0.012%
26	14.104	1830	1839	1851	rBV	3955320	5358696	55.08%	5.720%
27	14.393	1881	1888	1910	rBV	4521154	6419612	65.99%	6.852%
28	14.693	1931	1939	1949	rBV	2616662	3871377	39.80%	4.132%
29	14.904	1969	1975	1992	rBV	203554	438441	4.51%	0.468%
30	15.087	2002	2006	2010	rVB7	6623	10066	0.10%	0.011%
31	15.287	2034	2040	2045	rBV2	39136	59083	0.61%	0.063%
32	15.687	2099	2108	2121	rBV	5896857	8893630	91.42%	9.492%
33	15.804	2122	2128	2144	rVV	1603760	2274025	23.38%	2.427%
34	15.922	2144	2148	2160	rVB2	30600	65615	0.67%	0.070%
35	17.122	2349	2352	2357	rVB6	8123	11966	0.12%	0.013%
36	17.245	2370	2373	2376	rBV3	8502	12986	0.13%	0.014%

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37	17.457	2403	2409	2419	rBV	3317102	4741979	48.74%	5.061%
38	17.563	2420	2427	2450	rVB2	6125064	8882999	91.31%	9.481%
39	17.839	2472	2474	2478	rBV4	7171	10762	0.11%	0.011%
40	18.292	2547	2551	2557	rBV3	47922	81762	0.84%	0.087%
41	19.363	2730	2733	2738	rVB2	78101	94966	0.98%	0.101%
42	19.433	2741	2745	2749	rVV	67627	88467	0.91%	0.094%
43	19.528	2758	2761	2767	rBV6	46658	69656	0.72%	0.074%
44	19.792	2800	2806	2815	rBV	7641112	9728279	100.00%	10.383%
45	21.557	3101	3106	3114	rBV	3146715	4166521	42.83%	4.447%
46	23.257	3392	3395	3402	rVB	250774	375897	3.86%	0.401%
47	23.963	3506	3515	3528	rVB2	3488649	7885206	81.05%	8.416%
48	24.133	3538	3544	3554	rVB	1707194	3629340	37.31%	3.874%

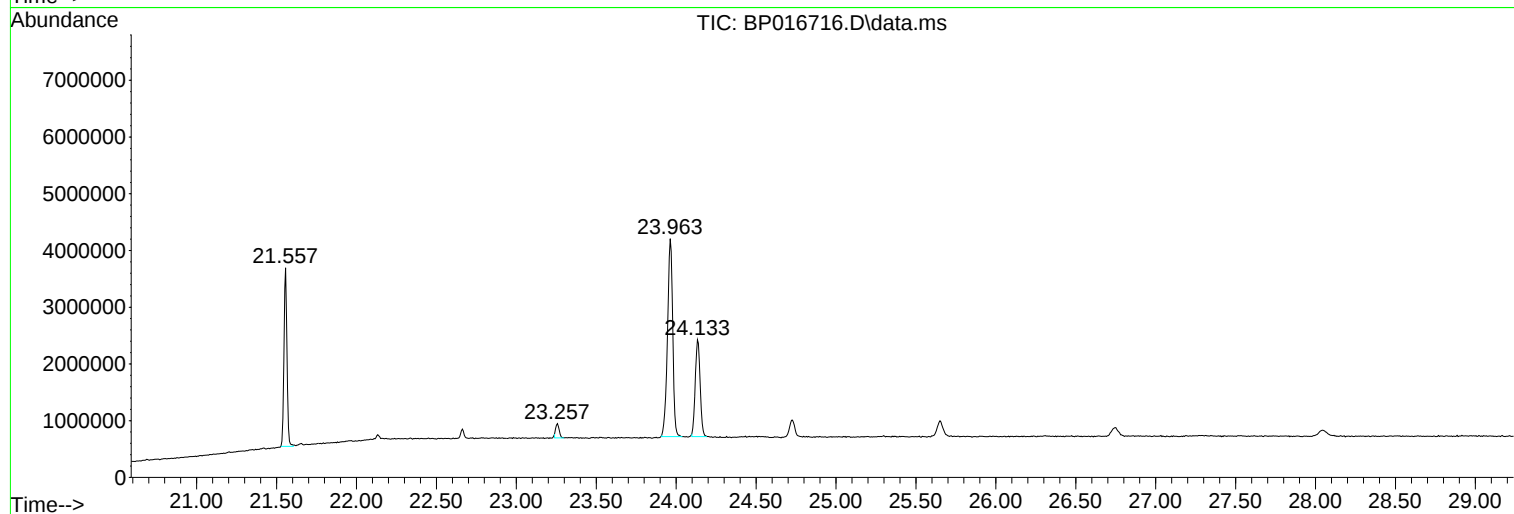
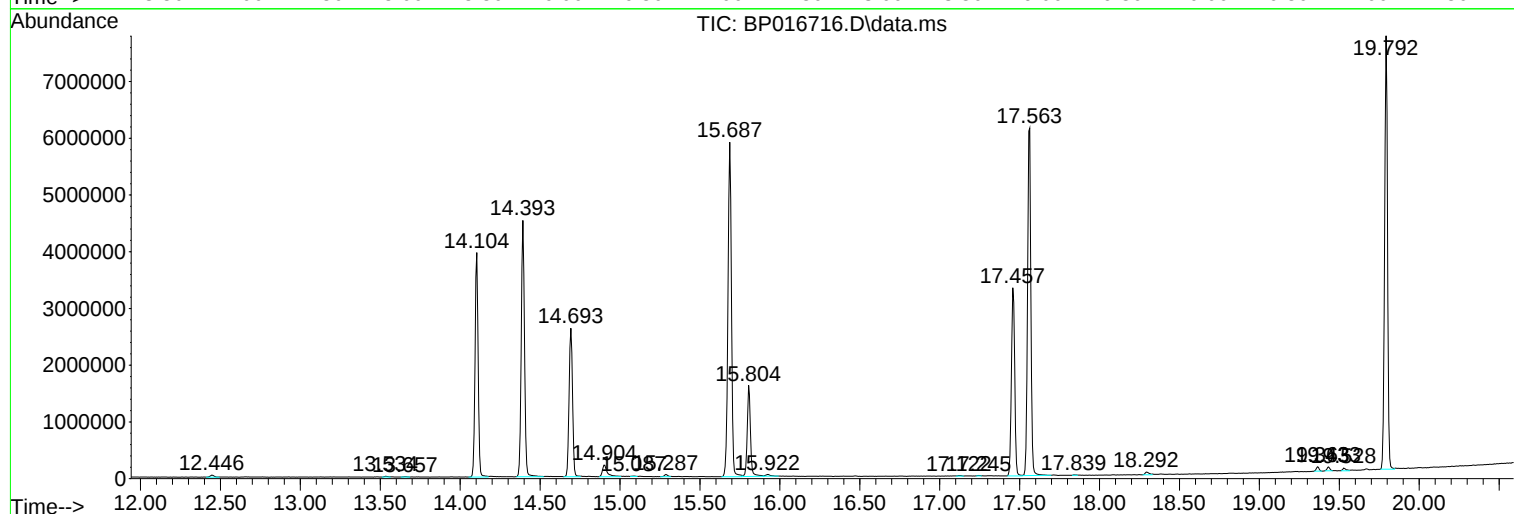
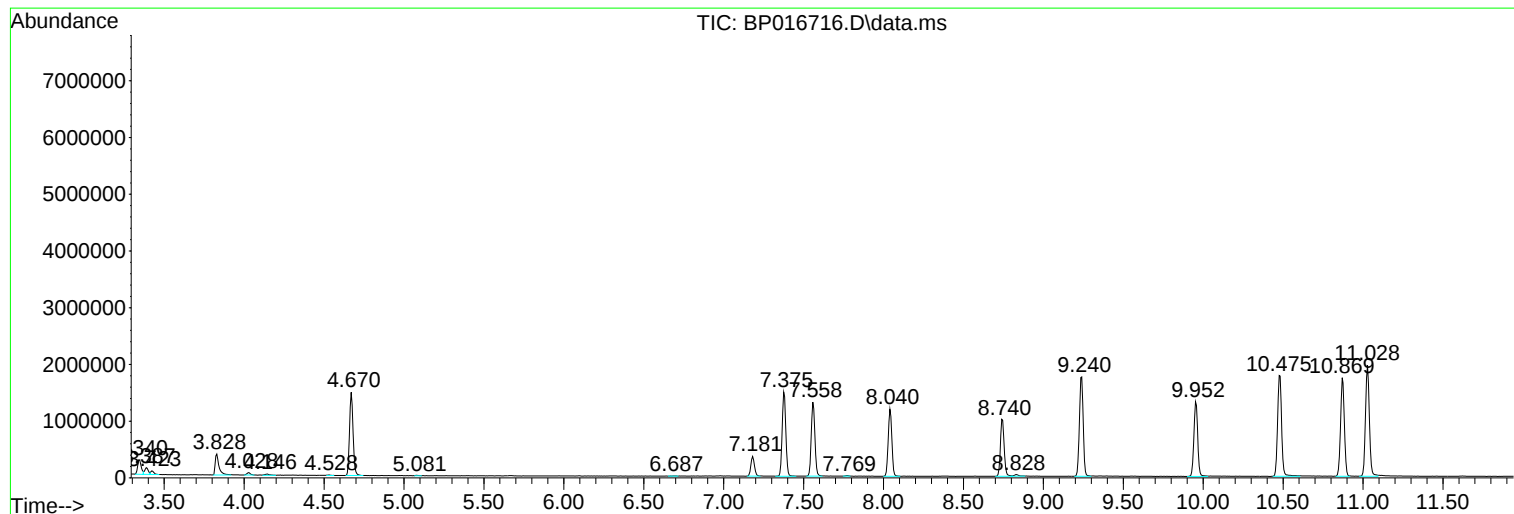
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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



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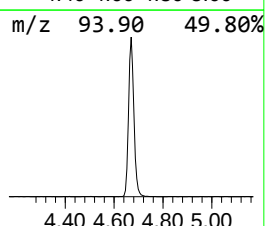
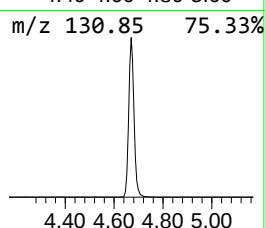
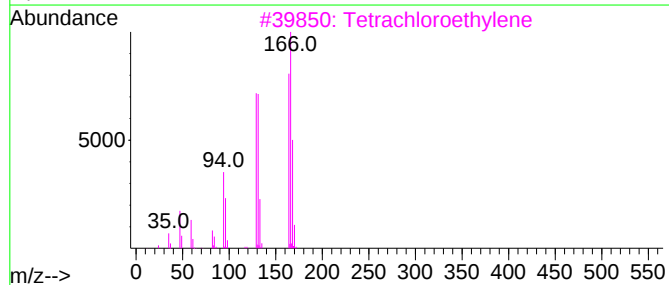
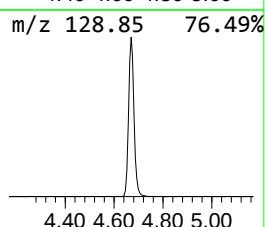
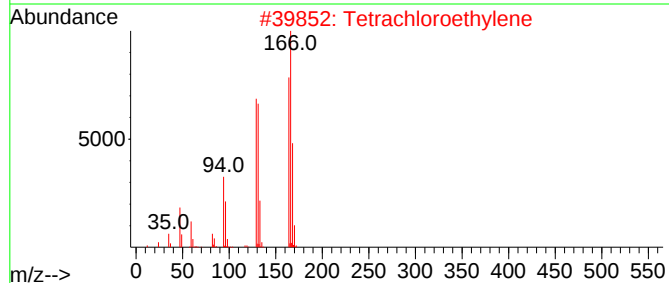
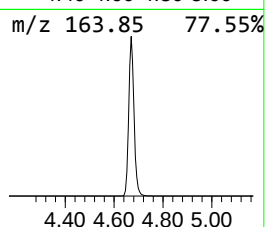
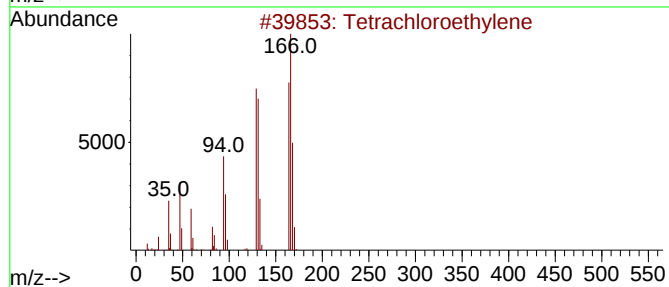
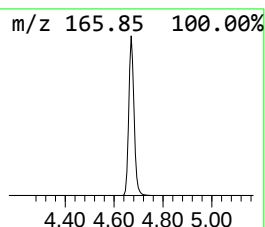
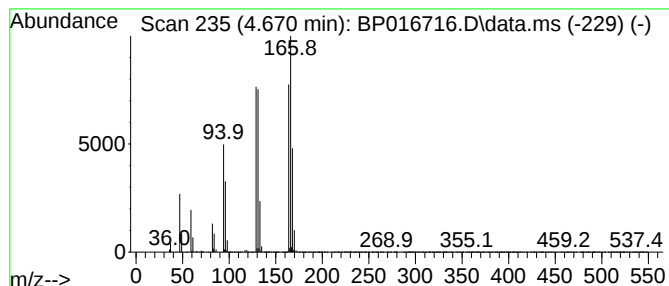
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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.
4.670	23.97 ng/ul	2223470	1,4-Dichlorobenzene-d4	8.040

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	98
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	38



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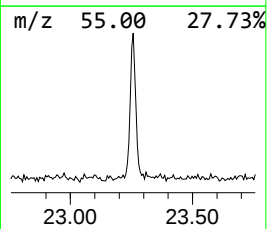
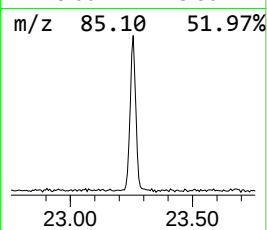
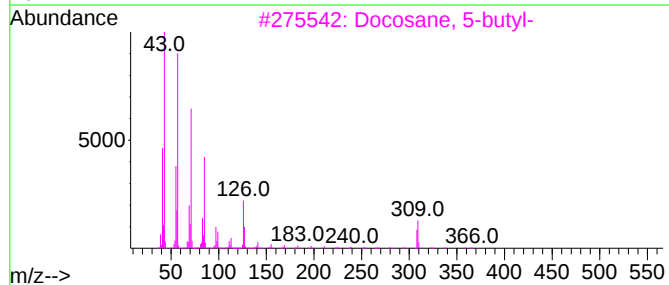
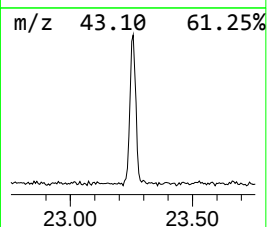
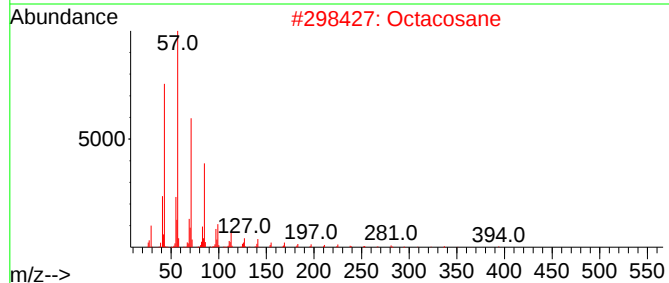
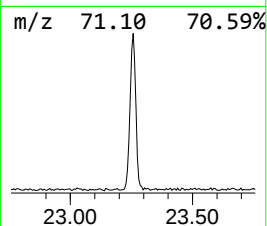
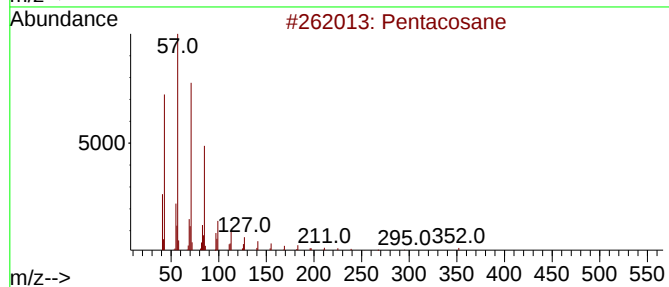
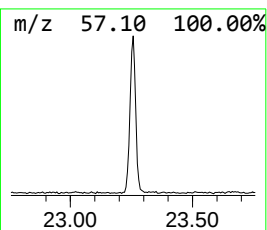
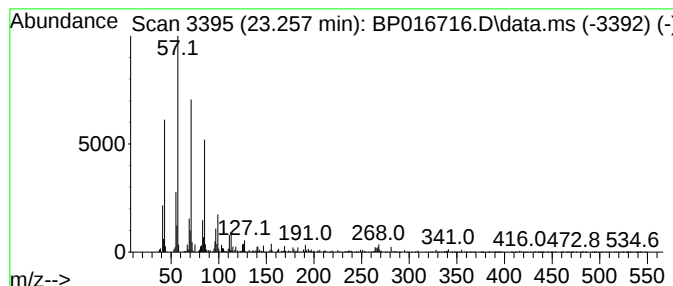
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.257	2.07 ng/ul	375897	Perylene-d12	24.133

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentacosane	352	C25H52	000629-99-2	93
2		Octacosane	394	C28H58	000630-02-4	93
3		Docosane, 5-butyl-	366	C26H54	055282-16-1	93
4		Docosane, 11-butyl-	366	C26H54	013475-76-8	93
5		Nonadecane	268	C19H40	000629-92-5	91



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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
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...	4.670	24.0	ng/u1	2223470	1	8.040	1855060	20.0
(DEL) Alkane: S...	23.257	2.1	ng/u1	375897	6	24.133	3629340	20.0