

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032023\
 Data File : BP014170.D
 Acq On : 21 Mar 2023 08:41
 Operator : CG/JU
 Sample : 01996-01
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0A24

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-BP030723.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014170.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.375	27	32	41	rBV2	235162	426559	2.48%	0.598%
2	3.446	41	44	53	rVB	76121	109785	0.64%	0.154%
3	3.852	111	113	130	rBV	136796	249279	1.45%	0.350%
4	4.716	253	260	274	rBV	11260307	17178610	100.00%	24.092%
5	7.210	677	684	697	rBV	142227	261570	1.52%	0.367%
6	7.369	704	711	722	rBV	632636	999856	5.82%	1.402%
7	7.575	740	746	756	rBV	570936	962072	5.60%	1.349%
8	8.046	819	826	835	rBV	607956	998021	5.81%	1.400%
9	8.763	941	948	967	rBV	462657	812759	4.73%	1.140%
10	9.222	1018	1026	1037	rBV	823906	1386190	8.07%	1.944%
11	9.604	1088	1091	1098	rVB4	12160	18642	0.11%	0.026%
12	9.946	1140	1149	1161	rBV	645228	1138816	6.63%	1.597%
13	10.493	1235	1242	1257	rBV	867885	1602115	9.33%	2.247%
14	10.869	1298	1306	1316	rBV	896012	1510051	8.79%	2.118%
15	11.010	1322	1330	1347	rBV	771261	1443685	8.40%	2.025%
16	12.457	1570	1576	1584	rBV2	15538	31721	0.18%	0.044%
17	13.575	1760	1766	1772	rBV	25201	40368	0.23%	0.057%
18	14.092	1846	1854	1869	rBV	2223490	3077323	17.91%	4.316%
19	14.387	1894	1904	1916	rBV	2289204	3478258	20.25%	4.878%
20	14.687	1949	1955	1964	rBV	1486825	2193343	12.77%	3.076%
21	14.916	1987	1994	2007	rBV3	74492	220089	1.28%	0.309%
22	15.681	2117	2124	2136	rBV	3138552	4512982	26.27%	6.329%
23	15.804	2138	2145	2156	rBV	1049401	1553194	9.04%	2.178%
24	15.922	2161	2165	2170	rVB2	18850	31473	0.18%	0.044%
25	16.045	2181	2186	2195	rVB	63839	92652	0.54%	0.130%
26	16.375	2237	2242	2251	rBV2	18730	35269	0.21%	0.049%
27	17.175	2374	2378	2383	rBV2	23602	30896	0.18%	0.043%
28	17.228	2384	2387	2394	rVB6	12563	22562	0.13%	0.032%
29	17.445	2417	2424	2434	rBV	1963093	2813533	16.38%	3.946%
30	17.545	2434	2441	2451	rBV	3653056	5170056	30.10%	7.251%
31	17.910	2500	2503	2508	rBV2	24908	31294	0.18%	0.044%
32	18.304	2566	2570	2580	rBV	154157	266480	1.55%	0.374%
33	18.410	2582	2588	2597	rVB	610891	857267	4.99%	1.202%
34	18.580	2614	2617	2622	rVB2	29430	33966	0.20%	0.048%
35	19.186	2717	2720	2723	rVB3	25458	28033	0.16%	0.039%
36	19.345	2743	2747	2751	rBV	55197	76195	0.44%	0.107%

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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

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37	19.410	2755	2758	2765	rVB	51552	81640	0.48%	0.114%
38	19.533	2775	2779	2789	rBV2	68875	125999	0.73%	0.177%
39	19.769	2813	2819	2825	rBV	4542217	6135460	35.72%	8.605%
40	21.192	3058	3061	3068	rBV2	236956	321622	1.87%	0.451%
41	21.521	3112	3117	3126	rBV	2300187	3141637	18.29%	4.406%
42	23.880	3511	3518	3530	rVB2	2584587	5171935	30.11%	7.253%
43	24.045	3539	3546	3555	rVB2	1193493	2631154	15.32%	3.690%

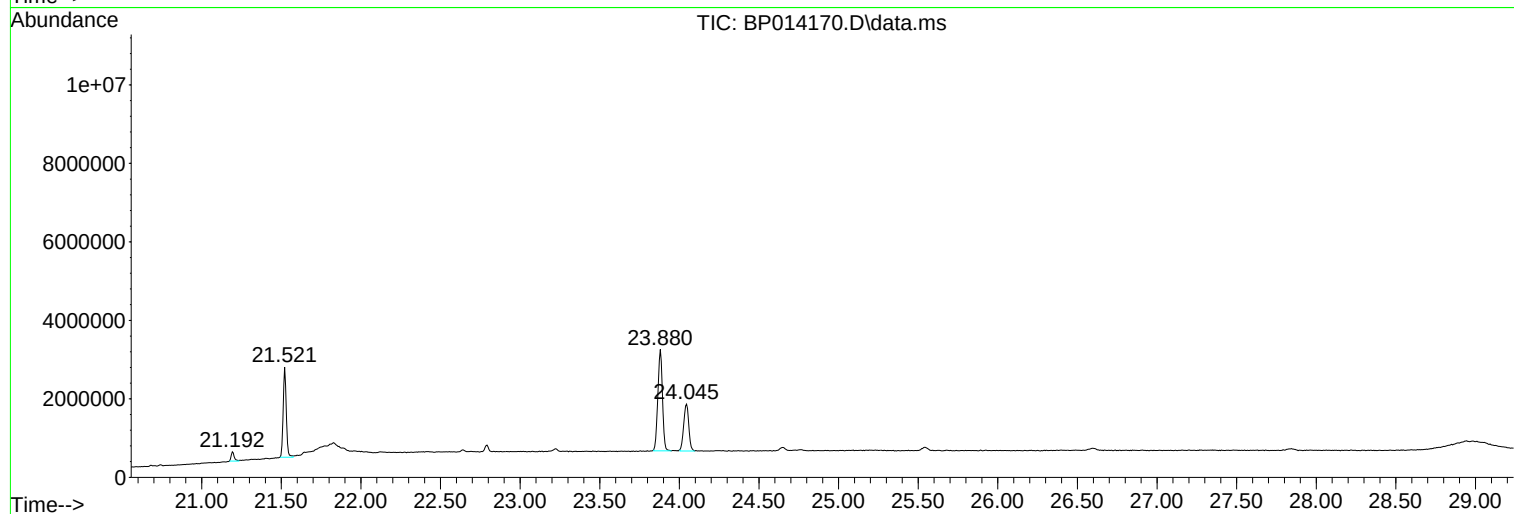
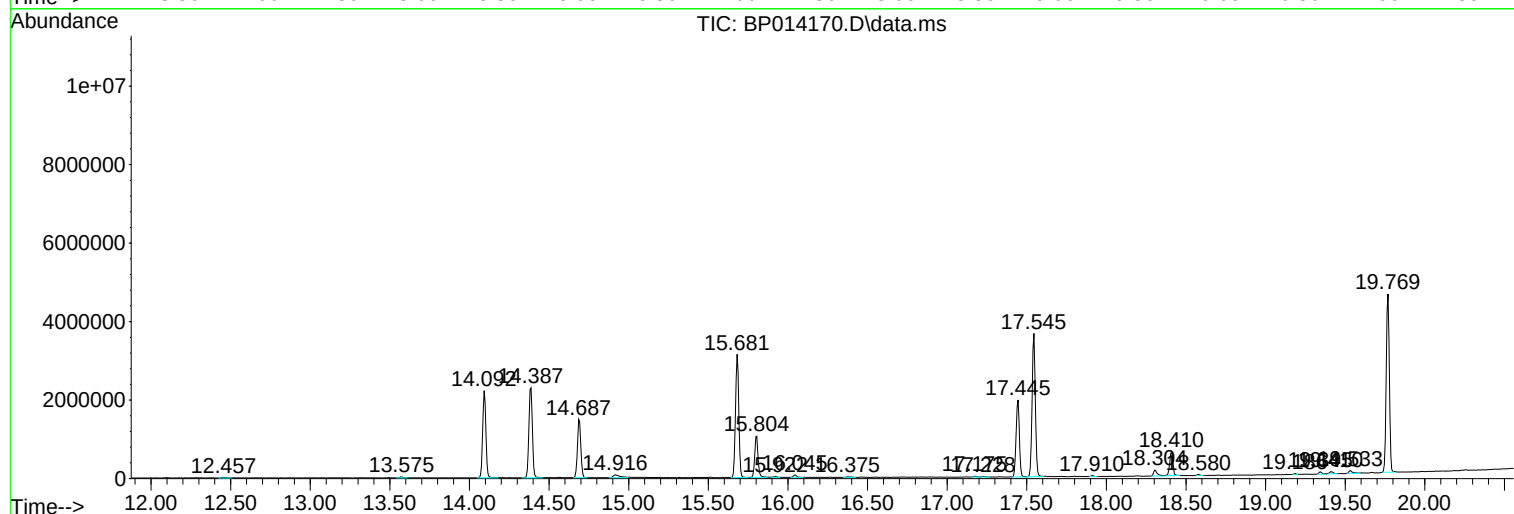
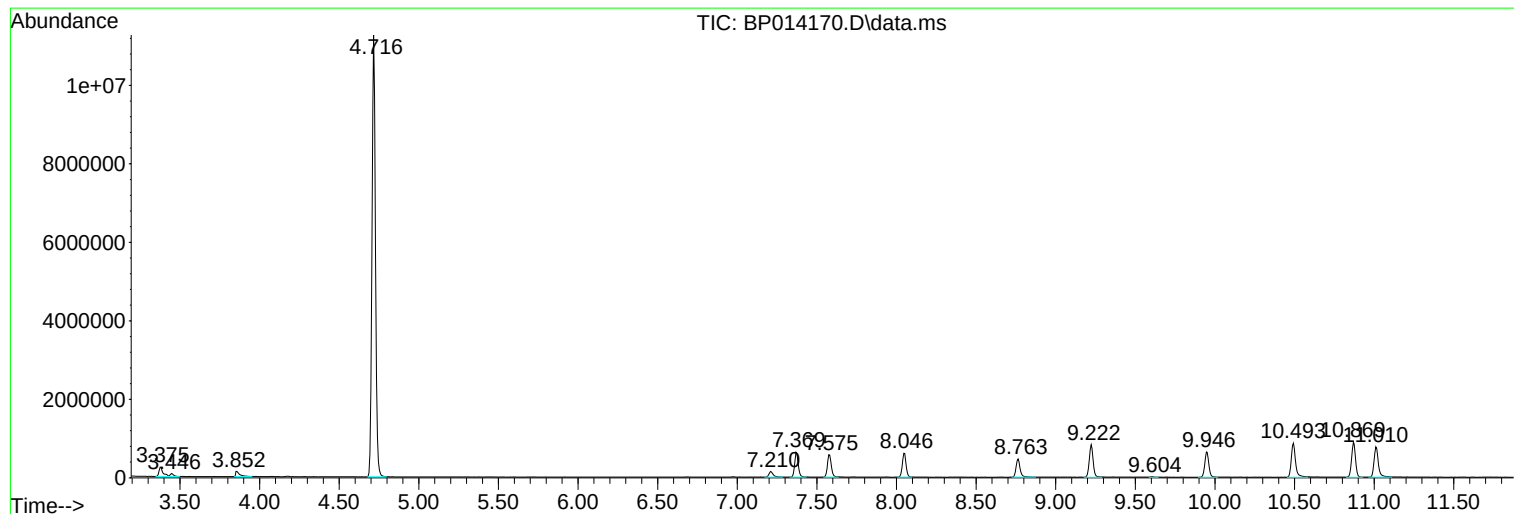
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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



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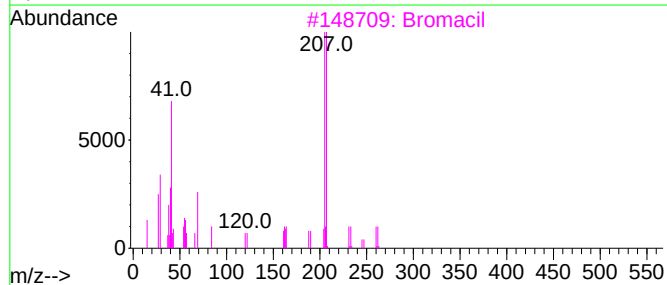
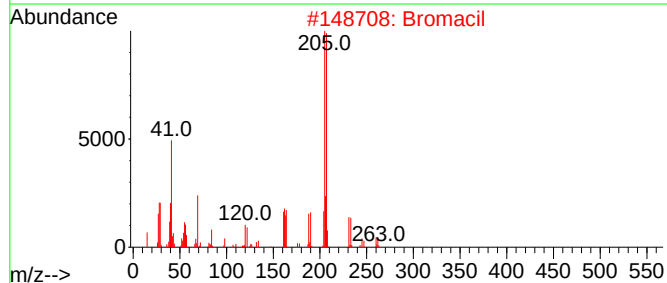
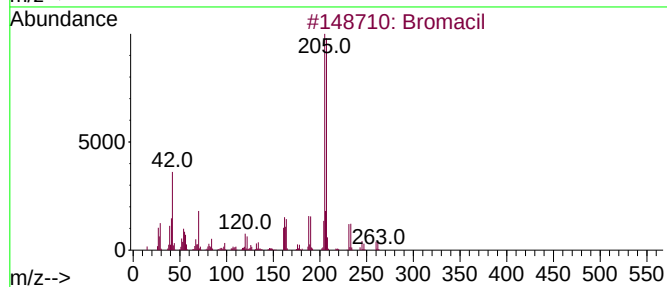
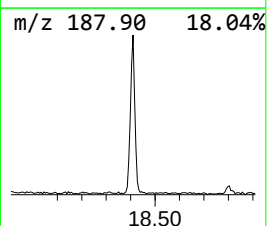
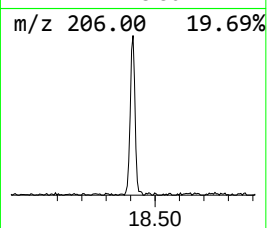
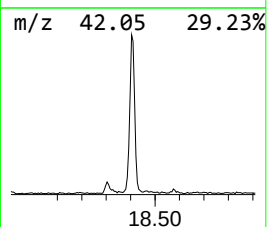
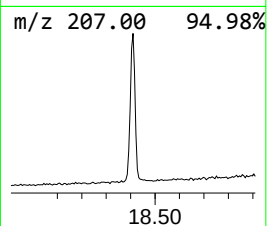
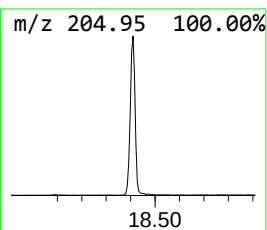
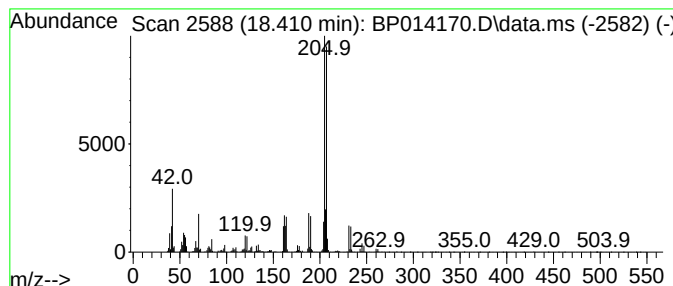
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Bromacil Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.410	6.09 ng/ul	857267	Phenanthrene-d10	17.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bromacil	260	C9H13BrN2O2	000314-40-9	94
2			Bromacil	260	C9H13BrN2O2	000314-40-9	91
3			Bromacil	260	C9H13BrN2O2	000314-40-9	87
4			5-Acetamido-2,4-dichlorobenzoic ...	247	C9H7Cl2NO3	050602-49-8	59
5			Benzoic acid, 3-amino-2,5-dichloro-	205	C7H5Cl2NO2	000133-90-4	50



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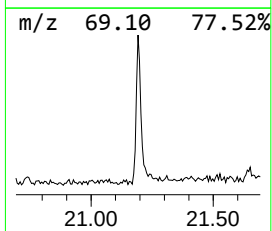
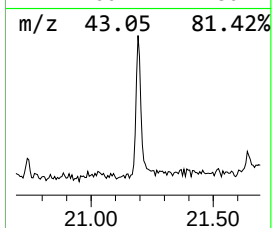
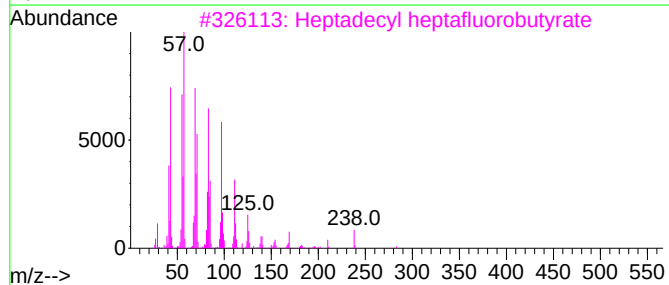
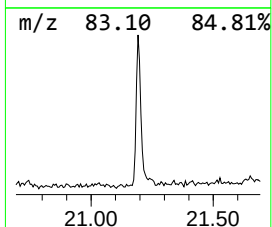
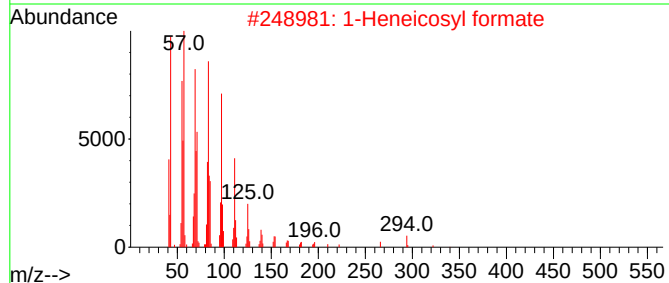
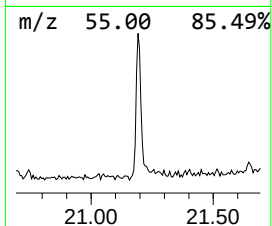
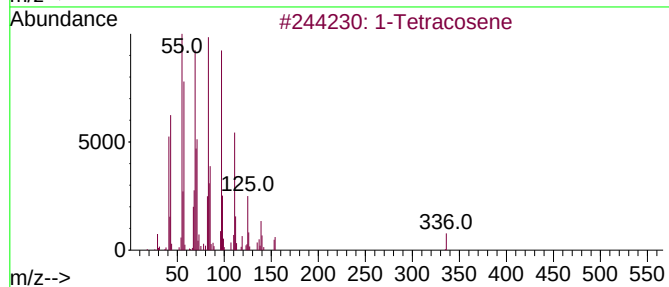
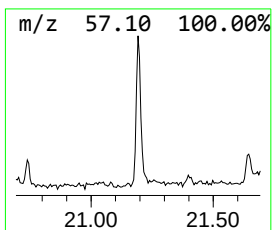
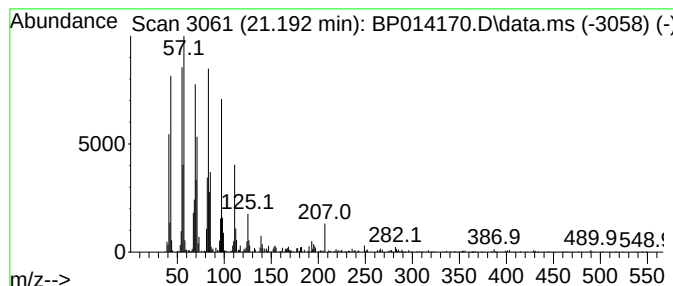
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Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.192	2.05 ng/ul	321622	Chrysene-d12		21.521	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Tetracosene		336	C24H48	010192-32-2	99
2	1-Heneicosyl formate		340	C22H44O2	077899-03-7	94
3	Heptadecyl heptafluorobutyrate		452	C21H35F7O2	959085-66-6	94
4	Pentafluoropropionic acid, penta...		374	C18H31F5O2	959092-08-1	94
5	Trifluoroacetoxy hexadecane		338	C18H33F3O2	006222-03-3	94



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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
Bromacil	18.410	6.1	ng/ul	857267	4	17.445	2813530	20.0
(DEL) Alkane: S...	21.192	2.0	ng/ul	321622	5	21.521	3141640	20.0