

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012056.D
 Acq On : 12 Oct 2022 00:18
 Operator : CG/JU
 Sample : N4991-05
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BF5

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-BP101122.M
 Title : SVOA CALIBRATION

Signal : TIC: BP012056.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.281	46	50	57	rVB	42231	56818	1.22%	0.126%
2	3.569	97	99	112	rVB	12356	20782	0.45%	0.046%
3	3.758	129	131	135	rBV3	9774	14178	0.30%	0.032%
4	3.805	136	139	143	rVB	17751	23891	0.51%	0.053%
5	3.928	155	160	172	rVB	50084	81372	1.75%	0.181%
6	4.022	172	176	182	rBV	25231	36887	0.79%	0.082%
7	4.399	237	240	247	rVB8	6883	9919	0.21%	0.022%
8	4.963	328	336	343	rBV2	34975	77645	1.67%	0.173%
9	5.487	421	425	432	rVB2	11150	22046	0.47%	0.049%
10	5.558	432	437	443	rVB4	4578	9603	0.21%	0.021%
11	5.858	480	488	494	rBV3	12982	24470	0.53%	0.054%
12	7.022	680	686	699	rBV	422544	696406	14.95%	1.548%
13	7.187	708	714	728	rBV	526853	861149	18.49%	1.914%
14	7.387	741	748	758	rBV	583592	946674	20.33%	2.104%
15	7.469	758	762	766	rBV	17917	26070	0.56%	0.058%
16	7.681	791	798	804	rVB8	3767	9511	0.20%	0.021%
17	7.857	817	828	842	rBV	765830	1320723	28.36%	2.936%
18	7.963	844	846	854	rVB9	4060	7714	0.17%	0.017%
19	8.569	943	949	961	rBV	385783	670516	14.40%	1.490%
20	8.693	966	970	975	rVB2	8994	13733	0.29%	0.031%
21	8.957	1011	1015	1019	rBV4	6209	10540	0.23%	0.023%
22	9.022	1019	1026	1033	rVV	544271	914320	19.63%	2.032%
23	9.075	1033	1035	1046	rVB	20545	33606	0.72%	0.075%
24	9.434	1087	1096	1100	rBV7	10049	20328	0.44%	0.045%
25	9.746	1141	1149	1164	rBV	411892	727681	15.63%	1.617%
26	10.287	1233	1241	1251	rBV	622860	1107011	23.77%	2.461%
27	10.446	1263	1268	1271	rBV2	7844	12938	0.28%	0.029%
28	10.663	1295	1305	1311	rBV	1029588	1774116	38.10%	3.943%
29	10.710	1311	1313	1319	rVB	33243	49553	1.06%	0.110%
30	10.810	1322	1330	1345	rBV	467750	915098	19.65%	2.034%
31	12.081	1541	1546	1550	rBV7	7252	10125	0.22%	0.023%
32	12.257	1571	1576	1582	rVB2	13146	23653	0.51%	0.053%
33	12.322	1582	1587	1595	rVB2	10992	21889	0.47%	0.049%
34	12.545	1621	1625	1631	rVB2	7488	12431	0.27%	0.028%
35	12.857	1673	1678	1682	rBV7	4349	7031	0.15%	0.016%
36	13.110	1716	1721	1725	rBV7	5602	9576	0.21%	0.021%

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Stop Thrs : 0

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
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37	13.322	1753	1757	1765	rVB6	9101	19850	0.43%	0.044%
38	13.828	1838	1843	1849	rBV9	11140	20467	0.44%	0.045%
39	13.916	1849	1858	1873	rBV	1149450	1643585	35.29%	3.653%
40	14.198	1899	1906	1926	rBV	1413250	2206410	47.38%	4.904%
41	14.392	1936	1939	1944	rVB3	7661	9407	0.20%	0.021%
42	14.504	1949	1958	1966	rBV	1502835	2211951	47.50%	4.916%
43	14.563	1966	1968	1972	rVB3	7570	10543	0.23%	0.023%
44	14.716	1988	1994	2005	rBV	132188	284428	6.11%	0.632%
45	14.904	2022	2026	2028	rBV	14495	21398	0.46%	0.048%
46	14.928	2028	2030	2037	rVB4	19804	29840	0.64%	0.066%
47	15.298	2089	2093	2095	rBV4	6330	9160	0.20%	0.020%
48	15.498	2119	2127	2134	rBV	2008263	2781299	59.73%	6.182%
49	15.551	2134	2136	2142	rVV3	17840	27462	0.59%	0.061%
50	15.622	2142	2148	2161	rVV	293844	431730	9.27%	0.960%
51	15.739	2165	2168	2177	rVB4	8490	17199	0.37%	0.038%
52	15.851	2183	2187	2192	rBV7	7304	13307	0.29%	0.030%
53	15.998	2209	2212	2219	rVB4	6475	10361	0.22%	0.023%
54	16.222	2245	2250	2257	rVB9	9077	23645	0.51%	0.053%
55	16.804	2343	2349	2353	rBV6	15448	27470	0.59%	0.061%
56	16.916	2364	2368	2375	rBV2	17240	29279	0.63%	0.065%
57	17.075	2393	2395	2403	rVB	13843	20774	0.45%	0.046%
58	17.263	2420	2427	2431	rBV	1804129	2652973	56.97%	5.897%
59	17.304	2431	2434	2438	rVV	344544	461287	9.91%	1.025%
60	17.363	2438	2444	2455	rVV	2017751	2982360	64.04%	6.629%
61	17.463	2456	2461	2470	rVB2	50032	86046	1.85%	0.191%
62	17.651	2490	2493	2494	rBV3	9376	10290	0.22%	0.023%
63	17.669	2494	2496	2507	rVB	21674	35505	0.76%	0.079%
64	17.928	2536	2540	2545	rBV	62130	77279	1.66%	0.172%
65	18.145	2571	2577	2580	rBV2	116327	195992	4.21%	0.436%
66	18.175	2580	2582	2587	rVB	56073	68910	1.48%	0.153%
67	18.316	2602	2606	2610	rBV2	42554	70857	1.52%	0.157%
68	18.616	2654	2657	2660	rVB	21425	25594	0.55%	0.057%
69	18.657	2661	2664	2668	rVB2	24181	29784	0.64%	0.066%
70	19.022	2722	2726	2731	rBV6	28618	59060	1.27%	0.131%
71	19.269	2764	2768	2774	rVB	543635	720821	15.48%	1.602%
72	19.374	2783	2786	2791	rBV2	79666	112861	2.42%	0.251%
73	19.598	2818	2824	2828	rBV	3006411	4331535	93.02%	9.628%
74	21.051	3067	3071	3080	rBV5	289569	512045	11.00%	1.138%
75	21.351	3116	3122	3126	rBV	2700967	3739292	80.30%	8.311%
76	23.610	3500	3506	3522	rVB2	2196039	4656716	100.00%	10.350%

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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

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Peak separation: 5

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77	23.768	3526	3533	3547	rVB2	1748827	3731937	80.14%	8.295%
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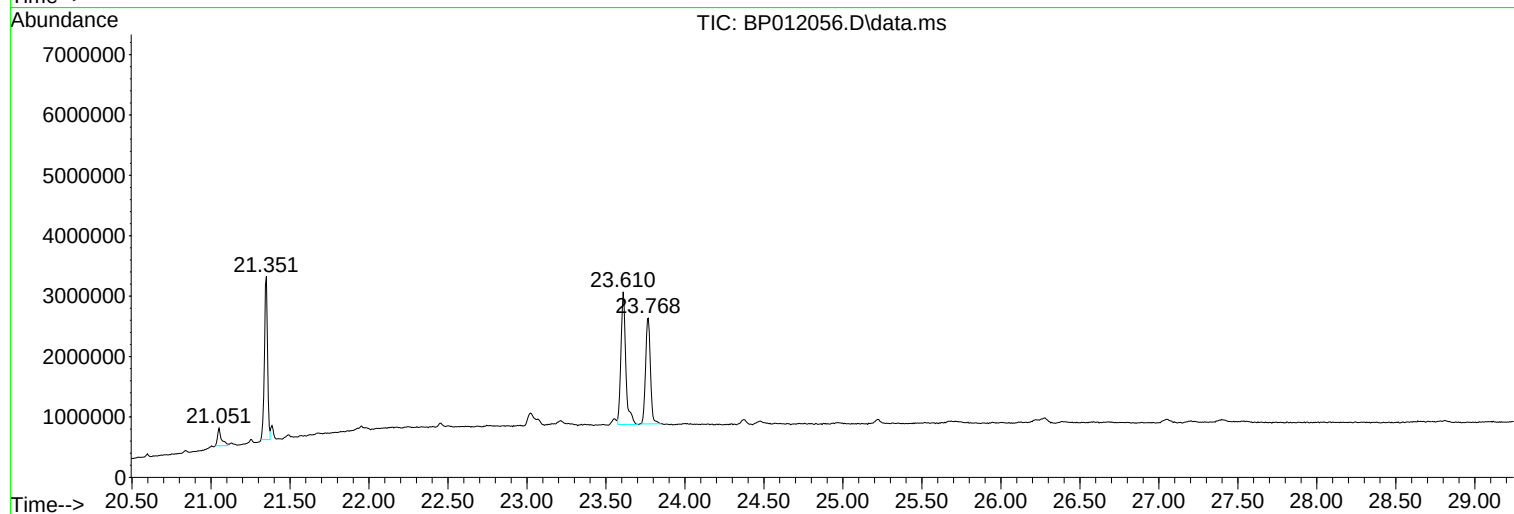
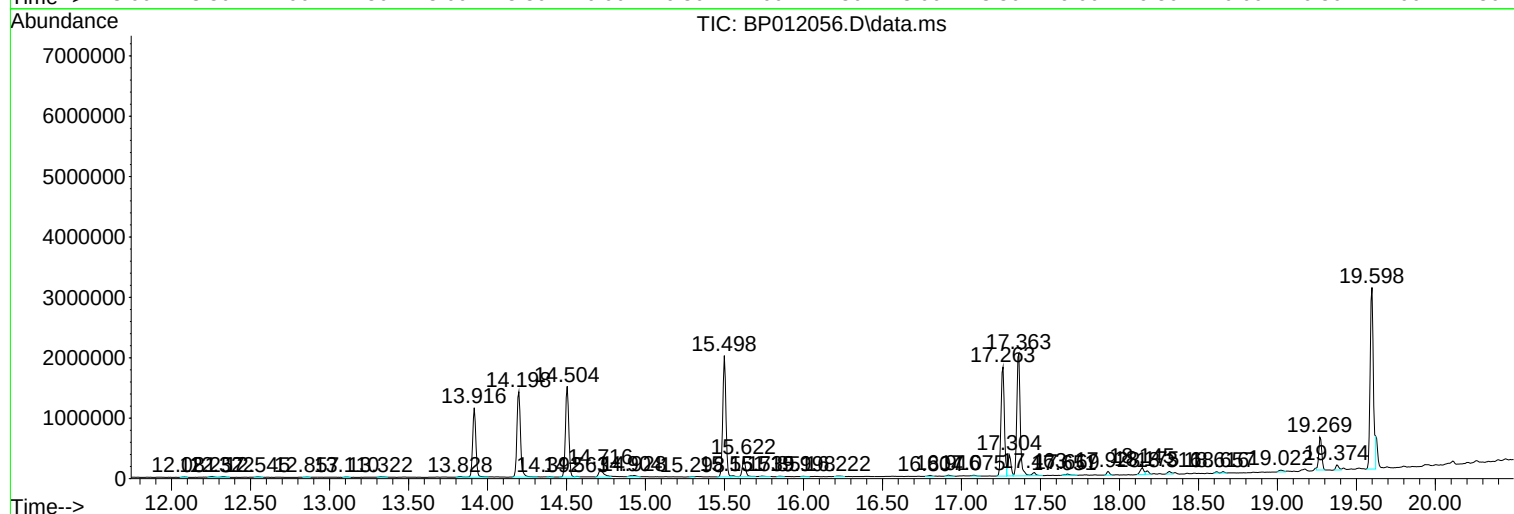
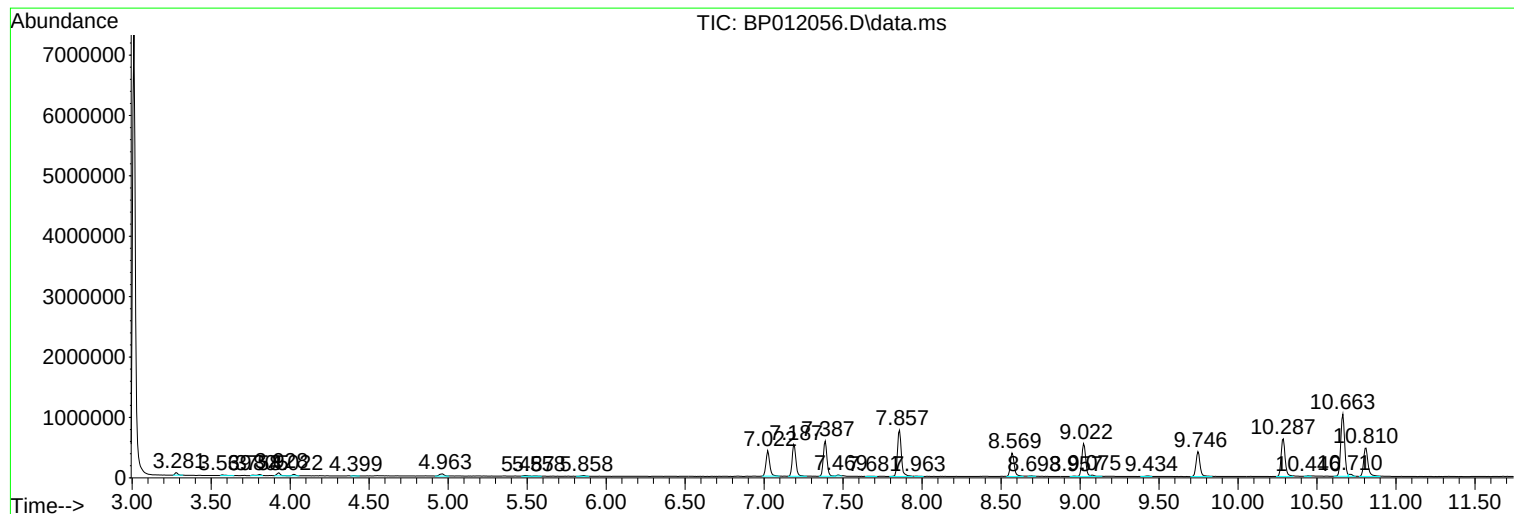
Sum of corrected areas: 44990712

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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



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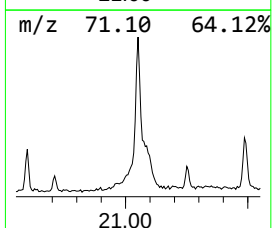
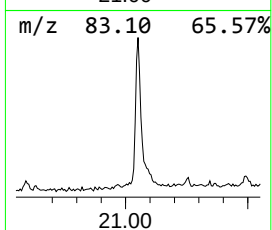
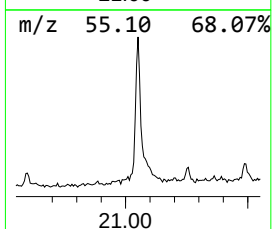
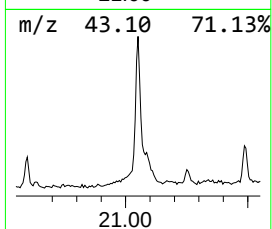
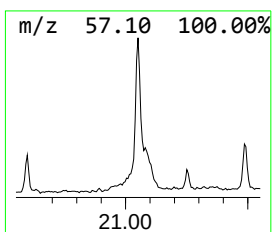
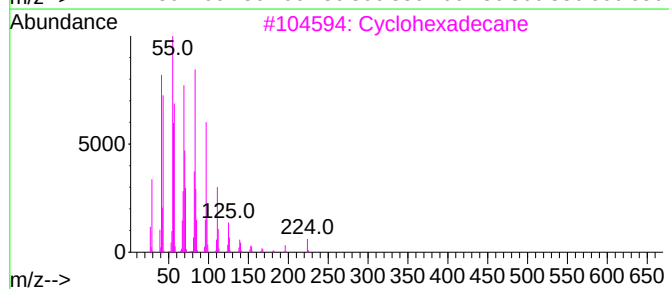
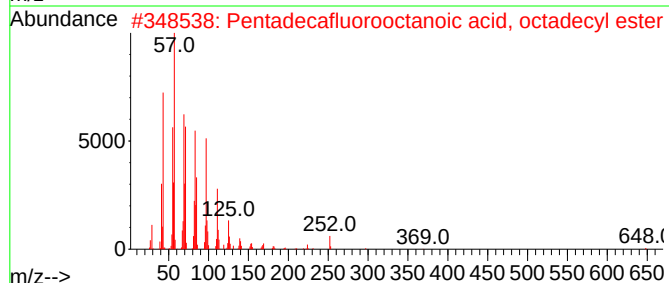
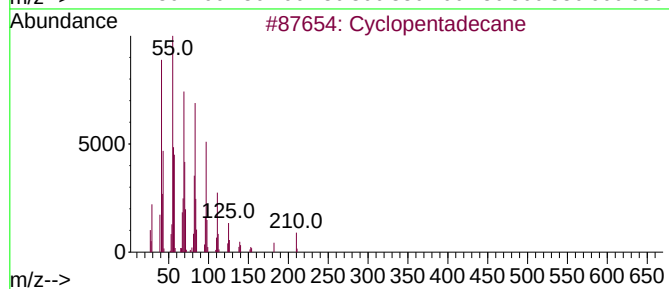
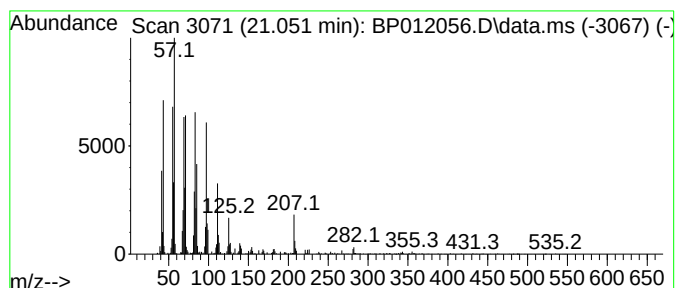
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Cyclic21.051 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.051	2.74 ng/ul	512045	Chrysene-d12	21.351	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentadecane	210	C15H30	000295-48-7	96
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
3	Cyclohexadecane	224	C16H32	000295-65-8	92
4	Eicosyl pentafluoropropionate	444	C23H41F5O2	1000351-80-8	90
5	Eicosane, 9-cyclohexyl-	364	C26H52	004443-61-2	90



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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
(DEL) Alkane: C...	21.051	2.7	ng/u1	512045	5	21.351	3739290	20.0