

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110221\
 Data File : BN017262.D
 Acq On : 03 Nov 2021 04:49
 Operator : CG/JU
 Sample : M4400-11
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BG346

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_N\Methods\SFAM-EPA-BN110221.M
 Title : SVOA CALIBRATION

Signal : TIC: BN017262.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.034	5	8	16	rVB	46513	63283	0.87%	0.101%
2	3.428	72	75	92	rBV	230451	422986	5.85%	0.672%
3	3.664	111	115	124	rVB	23643	37790	0.52%	0.060%
4	3.752	127	130	137	rVB	12483	20764	0.29%	0.033%
5	5.622	444	448	454	rBV	8231	11285	0.16%	0.018%
6	5.928	496	500	504	rBV4	5348	8217	0.11%	0.013%
7	6.699	623	631	642	rBV	186665	320373	4.43%	0.509%
8	6.858	652	658	673	rBV	708212	1117716	15.45%	1.776%
9	7.046	684	690	700	rBV	694373	1091321	15.08%	1.734%
10	7.511	762	769	777	rBV	784717	1206764	16.68%	1.917%
11	7.581	778	781	787	rVB2	9157	13761	0.19%	0.022%
12	8.228	884	891	901	rBV	562142	869021	12.01%	1.381%
13	8.663	958	965	974	rBV	849401	1363951	18.85%	2.167%
14	9.105	1035	1040	1049	rVB	17925	28257	0.39%	0.045%
15	9.375	1080	1086	1095	rBV	693000	1111394	15.36%	1.766%
16	9.546	1112	1115	1125	rVB4	5146	10293	0.14%	0.016%
17	9.910	1171	1177	1191	rBV	1028855	1679878	23.22%	2.669%
18	10.081	1200	1206	1222	rBV	99004	180672	2.50%	0.287%
19	10.281	1233	1240	1249	rBV	1205533	1972542	27.26%	3.134%
20	10.428	1256	1265	1282	rBV	1036062	1781159	24.62%	2.830%
21	11.887	1508	1513	1518	rBV	20680	32878	0.45%	0.052%
22	12.775	1660	1664	1669	rVB4	6042	8185	0.11%	0.013%
23	13.004	1697	1703	1708	rBV5	12969	21021	0.29%	0.033%
24	13.075	1711	1715	1719	rVB2	5251	7323	0.10%	0.012%
25	13.593	1796	1803	1815	rBV	2094098	2972334	41.08%	4.723%
26	13.851	1841	1847	1859	rBV	2451662	3623551	50.08%	5.758%
27	14.093	1881	1888	1894	rBV4	8315	14399	0.20%	0.023%
28	14.163	1894	1900	1908	rBV	1961850	2826426	39.06%	4.491%
29	14.334	1923	1929	1933	rBV6	7768	12368	0.17%	0.020%
30	14.393	1933	1939	1950	rBV2	98823	205201	2.84%	0.326%
31	14.551	1963	1966	1968	rBV3	5142	7439	0.10%	0.012%
32	14.610	1971	1976	1980	rVB4	4493	7577	0.10%	0.012%
33	15.051	2048	2051	2056	rBV4	10084	12362	0.17%	0.020%
34	15.169	2064	2071	2084	rBV	3235317	4642212	64.16%	7.376%
35	15.298	2086	2093	2102	rBV	1038672	1414293	19.55%	2.247%
36	15.404	2107	2111	2115	rVB2	22527	30400	0.42%	0.048%

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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_N\Methods\SFAM-EPA-BN110221.M
 Title : SVOA CALIBRATION

37	15.445	2115	2118	2120	rBV4	7087	7467	0.10%	0.012%
38	15.545	2129	2135	2140	rBV5	3979	7577	0.10%	0.012%
39	15.622	2140	2148	2155	rVB7	10553	19274	0.27%	0.031%
40	15.728	2162	2166	2171	rVB5	5929	10623	0.15%	0.017%

41	15.957	2201	2205	2211	rVB5	8128	13401	0.19%	0.021%
42	16.087	2221	2227	2229	rBV5	8692	15395	0.21%	0.024%
43	16.434	2283	2286	2290	rVB4	7086	8015	0.11%	0.013%
44	16.710	2326	2333	2336	rBV2	15667	27192	0.38%	0.043%
45	16.875	2355	2361	2363	rBV6	5935	8866	0.12%	0.014%

46	16.922	2363	2369	2378	rVV	2521619	3452202	47.71%	5.485%
47	17.022	2378	2386	2399	rVV	3696366	5309723	73.38%	8.437%
48	17.134	2402	2405	2410	rVV3	11543	19935	0.28%	0.032%
49	17.310	2431	2435	2440	rBV5	9764	17174	0.24%	0.027%
50	17.475	2458	2463	2467	rBV2	12681	20149	0.28%	0.032%

51	17.510	2467	2469	2475	rVB5	6724	8476	0.12%	0.013%
52	17.628	2482	2489	2494	rVB2	27871	37960	0.52%	0.060%
53	17.681	2494	2498	2504	rVB5	8948	14299	0.20%	0.023%
54	17.845	2521	2526	2530	rBV	63429	89291	1.23%	0.142%
55	17.916	2534	2538	2543	rVB	20144	27934	0.39%	0.044%

56	18.134	2570	2575	2577	rBV3	8280	13347	0.18%	0.021%
57	18.657	2659	2664	2666	rBV5	4676	8026	0.11%	0.013%
58	18.692	2666	2670	2673	rVV2	6308	7961	0.11%	0.013%
59	18.810	2687	2690	2693	rVB4	15753	16297	0.23%	0.026%
60	18.886	2699	2703	2710	rBV	49540	70628	0.98%	0.112%

61	18.957	2710	2715	2726	rBV	55621	90023	1.24%	0.143%
62	19.145	2743	2747	2753	rBV5	19260	32055	0.44%	0.051%
63	19.328	2772	2778	2789	rVV	4832411	6350043	87.76%	10.090%
64	20.416	2960	2963	2966	rVB2	18637	21146	0.29%	0.034%
65	20.880	3037	3042	3057	rBV	193609	469636	6.49%	0.746%

66	21.075	3072	3075	3078	rBV	21020	22789	0.31%	0.036%
67	21.133	3079	3085	3097	rVV	3174674	4072040	56.28%	6.470%
68	21.322	3114	3117	3120	rBV2	29407	27662	0.38%	0.044%
69	21.751	3187	3190	3198	rBV2	49958	76146	1.05%	0.121%
70	22.204	3264	3267	3277	rVB	77569	106354	1.47%	0.169%

71	22.710	3349	3353	3361	rVB	142338	198202	2.74%	0.315%
72	23.192	3427	3435	3443	rBV	4057221	7235915	100.00%	11.497%
73	23.269	3444	3448	3452	rVV	184761	310860	4.30%	0.494%
74	23.333	3452	3459	3472	rVB2	2520698	4608830	63.69%	7.323%
75	23.910	3551	3557	3566	rBV	182382	354096	4.89%	0.563%

76	24.651	3678	3683	3693	rVB2	146640	304971	4.21%	0.485%
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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

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : SVOA CALIBRATION

77	25.521	3825	3831	3841	rVB2	104120	274637	3.80%	0.436%
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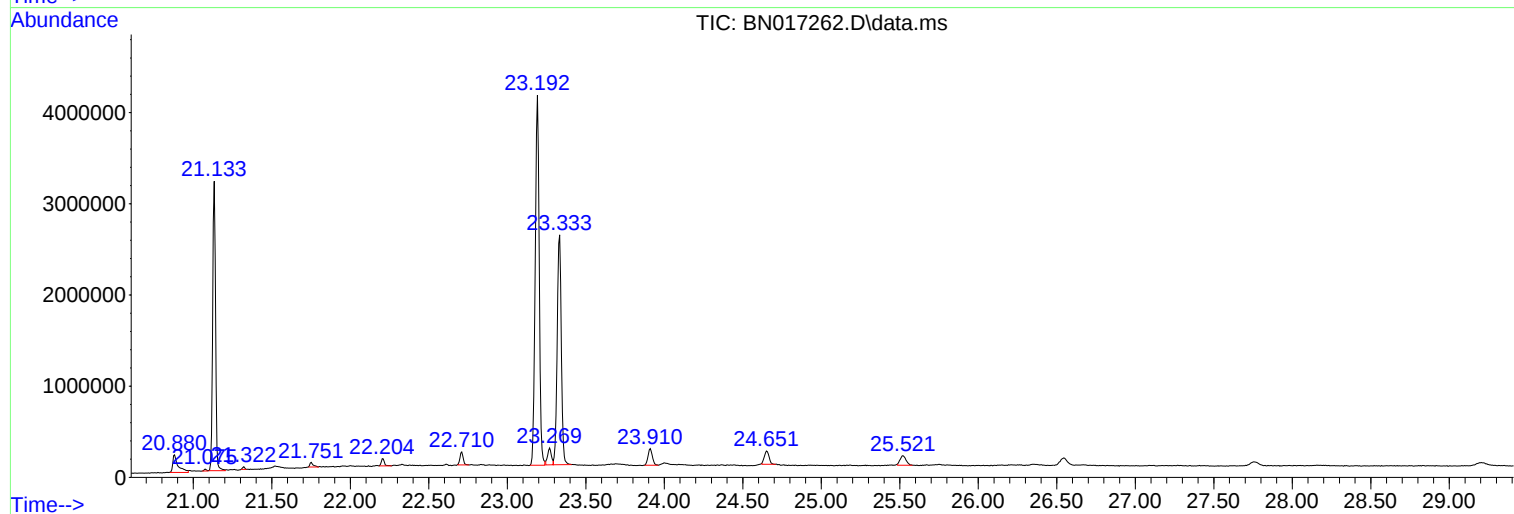
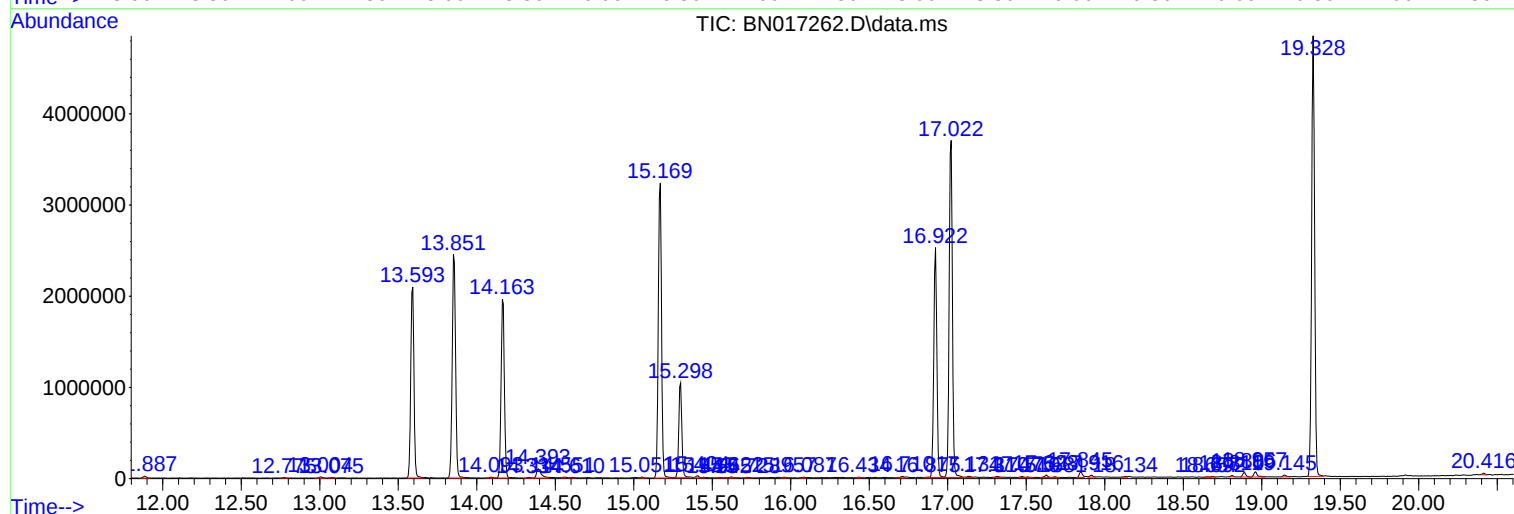
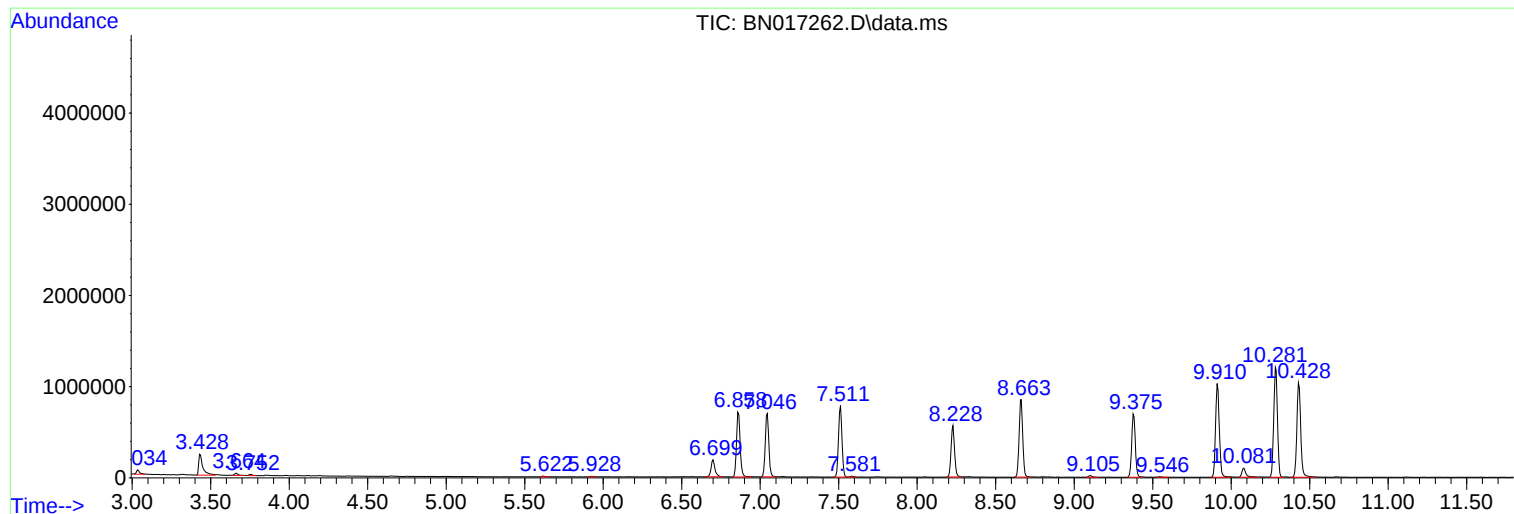
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M
Quant Title : SVOA CALIBRATION

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



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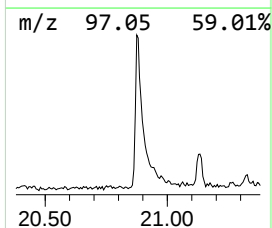
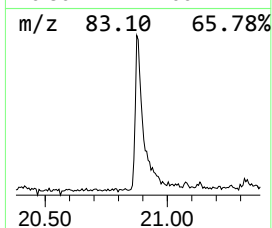
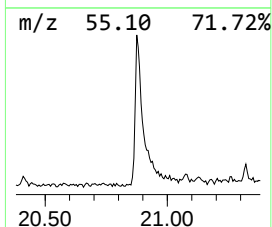
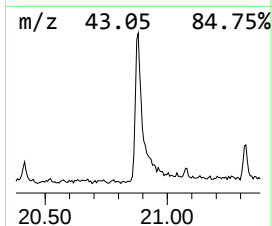
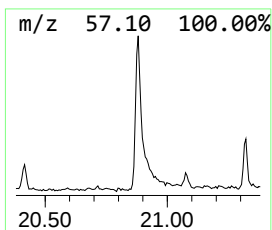
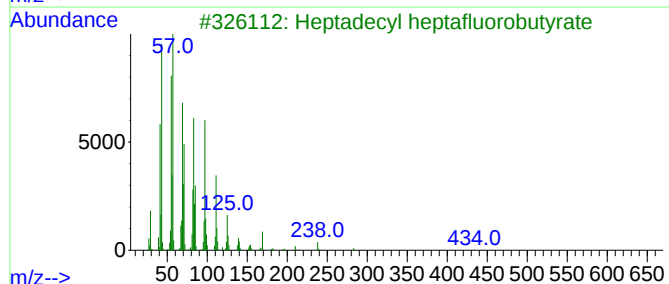
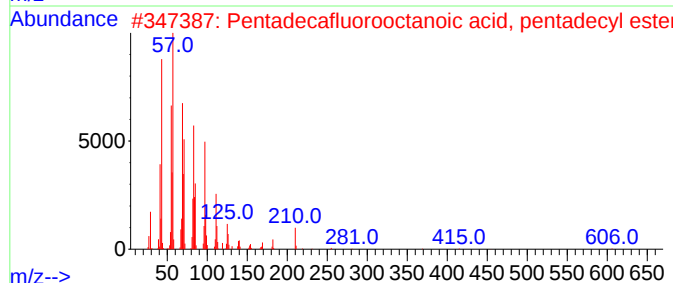
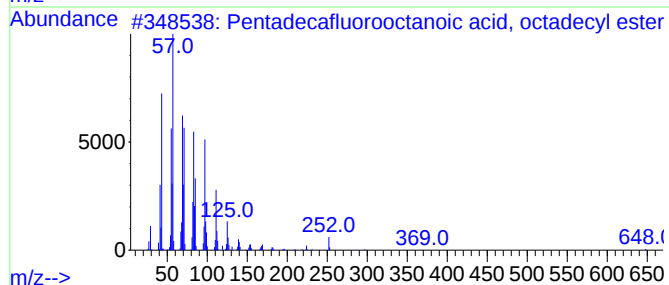
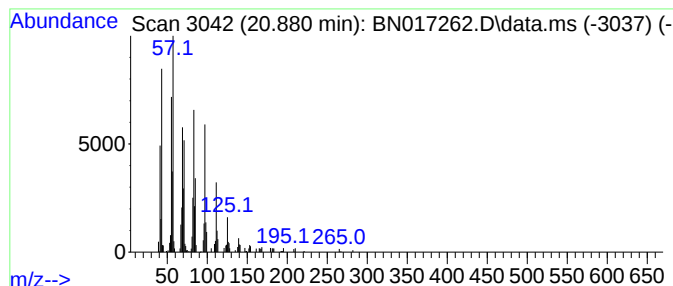
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN110221.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Pentadecafluorooctanoic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.880	2.31 ng/ul	469636	Chrysene-d12	21.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	95
2			Pentadecafluorooctanoic acid, pe...	624	C23H31F15O2	1000406-04-5	94
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
4			1-Hexacosene	364	C26H52	018835-33-1	93
5			1-Heneicosanol	312	C21H44O	015594-90-8	93



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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
Pentadecafluoro...	20.880	2.3	ng/u1	469636	5	21.133	4072040	20.0