

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012066.D
 Acq On : 12 Oct 2022 06:34
 Operator : CG/JU
 Sample : N4991-11
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 COBL4

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: BP012066.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.275	46	49	56	rVB	41754	54861	1.35%	0.132%
2	3.581	99	101	109	rVB7	4824	8747	0.21%	0.021%
3	3.805	134	139	148	rVB4	14362	26222	0.64%	0.063%
4	3.922	155	159	165	rVB	45832	69198	1.70%	0.166%
5	4.022	172	176	182	rVB	29014	39430	0.97%	0.095%
6	4.399	238	240	245	rVB6	4928	6957	0.17%	0.017%
7	4.958	328	335	343	rBV2	34311	72609	1.78%	0.174%
8	5.058	347	352	358	rVB9	3339	7179	0.18%	0.017%
9	5.352	398	402	407	rVB3	7039	9639	0.24%	0.023%
10	5.481	419	424	432	rBV	14265	32186	0.79%	0.077%
11	5.852	480	487	496	rVB	13908	26835	0.66%	0.064%
12	6.822	648	652	658	rVB4	4701	8637	0.21%	0.021%
13	6.940	667	672	675	rBV5	6067	10130	0.25%	0.024%
14	7.022	678	686	699	rVV	322760	551405	13.54%	1.323%
15	7.187	707	714	727	rBV	518267	845954	20.77%	2.029%
16	7.381	740	747	758	rBV	501331	817079	20.06%	1.960%
17	7.469	758	762	764	rVV2	12466	20707	0.51%	0.050%
18	7.499	764	767	773	rVB2	14424	21111	0.52%	0.051%
19	7.852	818	827	842	rBV	808696	1360728	33.41%	3.264%
20	7.958	842	845	852	rVB9	6598	15153	0.37%	0.036%
21	8.569	942	949	961	rBV	317948	561846	13.79%	1.348%
22	8.687	965	969	976	rVB3	6704	13332	0.33%	0.032%
23	8.963	1010	1016	1019	rBV6	7319	13073	0.32%	0.031%
24	9.022	1019	1026	1046	rVB	522081	919388	22.57%	2.205%
25	9.422	1087	1094	1101	rBV9	8803	17391	0.43%	0.042%
26	9.746	1141	1149	1167	rBV	383655	682645	16.76%	1.637%
27	10.281	1233	1240	1254	rBV	545780	976394	23.97%	2.342%
28	10.440	1262	1267	1271	rBV5	8054	12706	0.31%	0.030%
29	10.557	1284	1287	1293	rBV5	4452	7416	0.18%	0.018%
30	10.657	1295	1304	1311	rBV	1089327	1888223	46.36%	4.529%
31	10.804	1322	1329	1349	rBV	420599	798234	19.60%	1.915%
32	12.075	1542	1545	1551	rBV6	4390	6833	0.17%	0.016%
33	12.251	1564	1575	1583	rVB	13305	29077	0.71%	0.070%
34	12.328	1583	1588	1595	rBV2	7463	15809	0.39%	0.038%
35	12.546	1619	1625	1632	rVB3	5219	9997	0.25%	0.024%
36	13.104	1716	1720	1724	rVB7	4902	7417	0.18%	0.018%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
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37	13.316	1751	1756	1764	rBV8	5329	11363	0.28%	0.027%
38	13.822	1835	1842	1847	rBV8	8190	17941	0.44%	0.043%
39	13.910	1851	1857	1872	rVV	1087821	1601944	39.33%	3.842%
40	14.193	1898	1905	1918	rBV	1381976	2096141	51.46%	5.028%
41	14.504	1948	1958	1971	rBV	1578931	2389127	58.66%	5.730%
42	14.716	1988	1994	2012	rBV	58834	172960	4.25%	0.415%
43	14.904	2022	2026	2027	rBV	7118	8266	0.20%	0.020%
44	14.928	2027	2030	2036	rVB5	15270	22890	0.56%	0.055%
45	15.498	2120	2127	2141	rBV	1913930	2779550	68.24%	6.667%
46	15.622	2142	2148	2162	rVV	222417	348348	8.55%	0.836%
47	15.739	2165	2168	2174	rVB2	8632	13802	0.34%	0.033%
48	16.210	2244	2248	2257	rVB2	6982	13369	0.33%	0.032%
49	16.363	2268	2274	2277	rBV8	3800	8546	0.21%	0.020%
50	16.798	2344	2348	2352	rBV4	10851	15072	0.37%	0.036%
51	16.916	2364	2368	2370	rBV4	6539	8638	0.21%	0.021%
52	17.257	2420	2426	2432	rBV	1895462	2807252	68.92%	6.733%
53	17.298	2432	2433	2437	rVV	57019	61476	1.51%	0.147%
54	17.357	2437	2443	2456	rVV	2063449	2976211	73.07%	7.139%
55	17.457	2456	2460	2469	rVB2	47280	71316	1.75%	0.171%
56	17.651	2489	2493	2497	rBV7	6009	13094	0.32%	0.031%
57	17.928	2536	2540	2544	rBV3	16404	22046	0.54%	0.053%
58	18.139	2572	2576	2586	rBV2	99278	174068	4.27%	0.418%
59	19.169	2749	2751	2757	rVB2	25789	33939	0.83%	0.081%
60	19.239	2760	2763	2765	rBV	34247	44178	1.08%	0.106%
61	19.269	2765	2768	2773	rVB	85685	114824	2.82%	0.275%
62	19.375	2783	2786	2792	rBV3	48914	73826	1.81%	0.177%
63	19.592	2818	2823	2835	rBV	2879619	4073112	100.00%	9.770%
64	21.051	3067	3071	3080	rBV2	311669	498153	12.23%	1.195%
65	21.345	3116	3121	3126	rBV	2715689	3535601	86.80%	8.480%
66	23.610	3500	3506	3520	rVB2	2016978	4034285	99.05%	9.676%
67	23.763	3526	3532	3543	rVB2	1855127	3686129	90.50%	8.841%

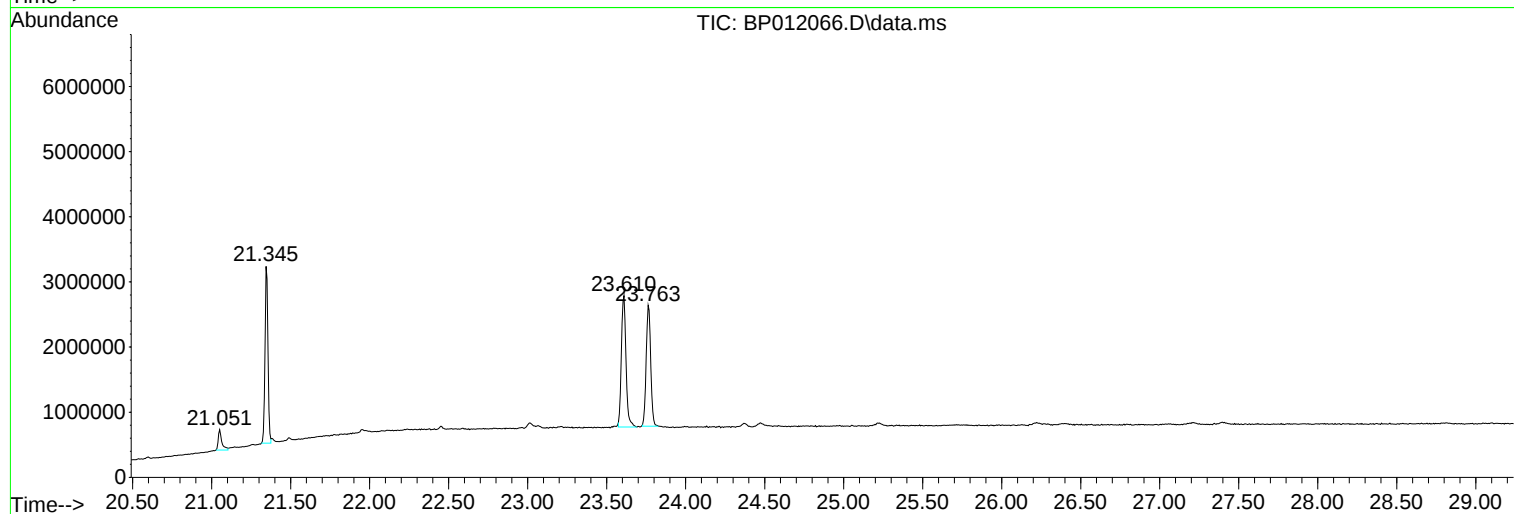
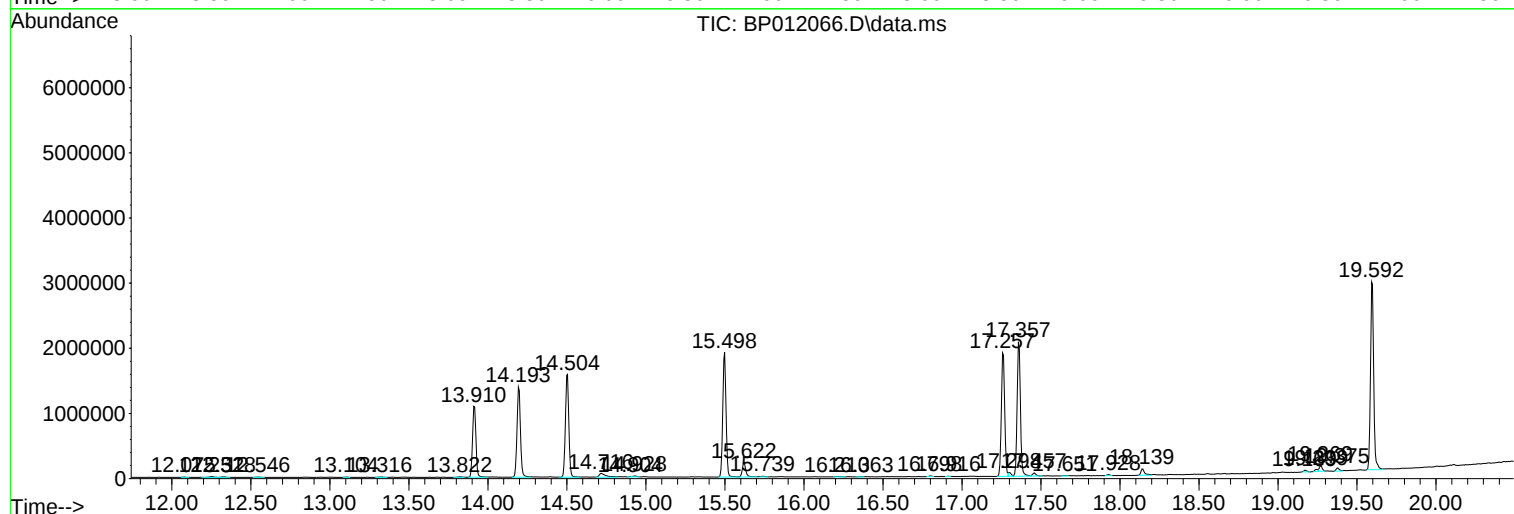
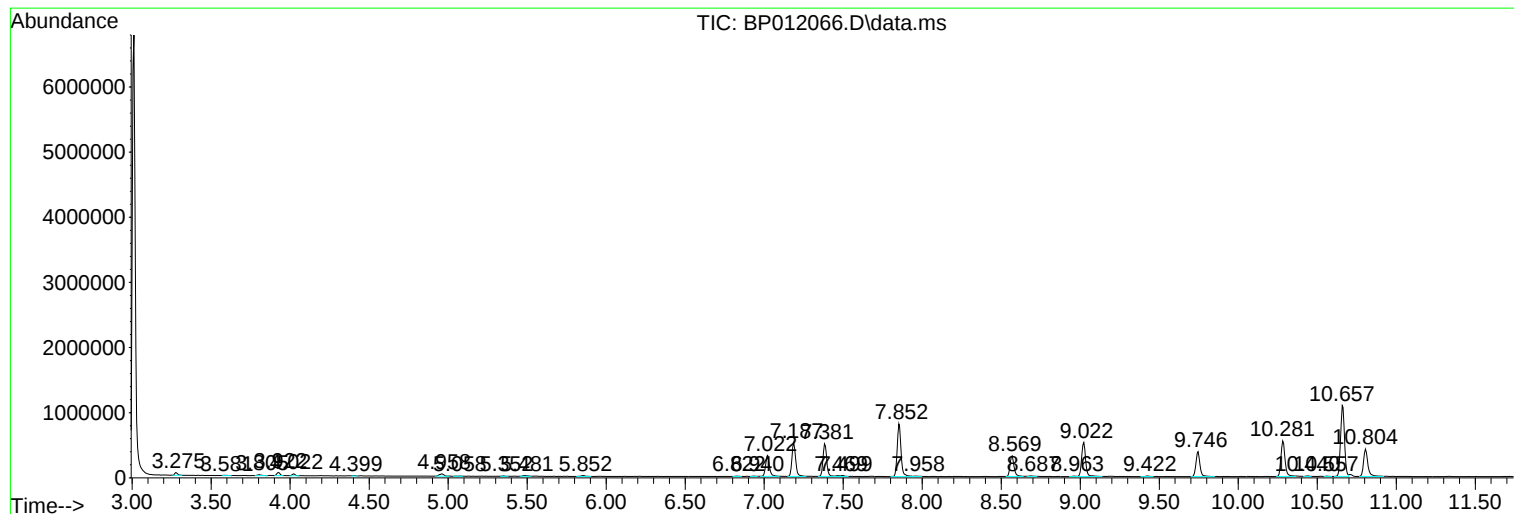
Sum of corrected areas: 41692015

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ClientSampleId :
C0BL4

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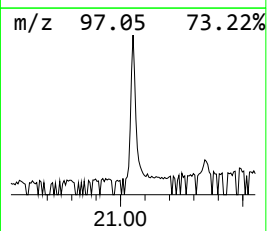
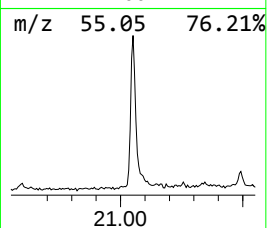
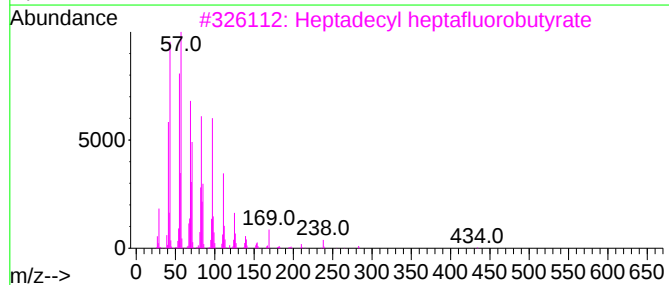
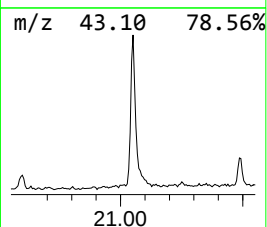
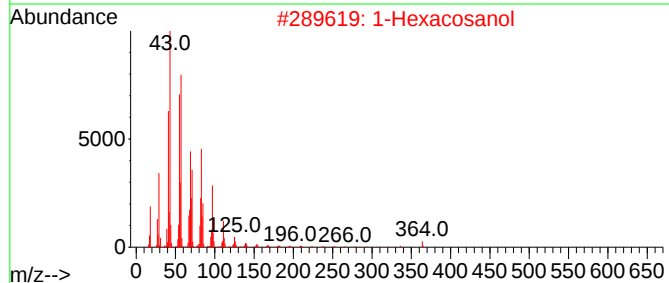
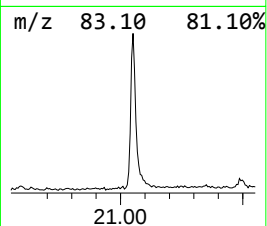
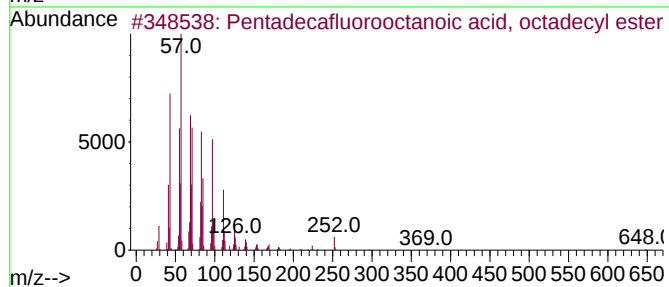
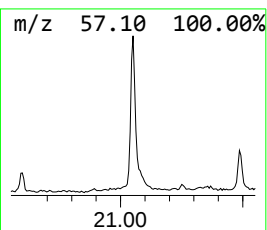
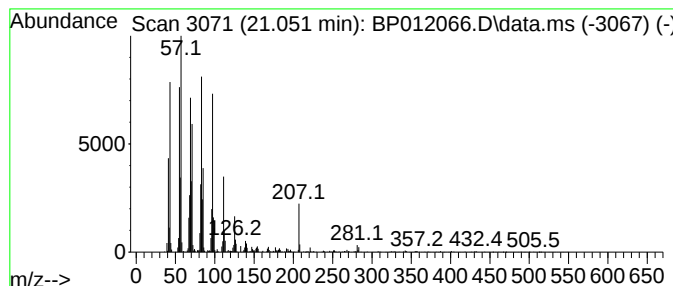
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Pentadecafluorooctanoic aci... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.051	2.82 ng/ul	498153	Chrysene-d12	21.345	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93
2	1-Hexacosanol	382	C26H54O	000506-52-5	90
3	Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	90
4	1-Heptacosanol	396	C27H56O	002004-39-9	90
5	Nonadecyl pentafluoropropionate	430	C22H39F5O2	1000351-88-8	90



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Pentadecafluoro...	21.051	2.8	ng/u1	498153	5	21.345	3535600	20.0