

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP072122\
 Data File : BP011071.D
 Acq On : 21 Jul 2022 20:13
 Operator : CG/JU
 Sample : N3709-03
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 H0B19

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-BP071422.M
 Title : SVOA CALIBRATION

Signal : TIC: BP011071.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	12	16	23	rVB	100272	144847	2.18%	0.187%
2	3.452	60	62	72	rVB	57769	86797	1.30%	0.112%
3	3.593	83	86	99	rBV	358007	569832	8.56%	0.737%
4	3.687	99	102	109	rVB2	57290	85904	1.29%	0.111%
5	3.805	118	122	129	rVB	93107	137689	2.07%	0.178%
6	3.899	133	138	144	rVB	210060	282132	4.24%	0.365%
7	4.446	228	231	235	rBV6	14038	19957	0.30%	0.026%
8	4.822	288	295	299	rBV2	91490	178588	2.68%	0.231%
9	5.275	369	372	378	rVB2	9799	13239	0.20%	0.017%
10	5.399	389	393	399	rVB	25034	36218	0.54%	0.047%
11	5.781	455	458	464	rVB7	6476	12441	0.19%	0.016%
12	5.934	480	484	487	rBV6	6333	8395	0.13%	0.011%
13	6.851	631	640	655	rBV	1340466	2092320	31.44%	2.707%
14	7.016	661	668	676	rVV	1141872	1738745	26.12%	2.250%
15	7.093	677	681	688	rVB	26019	45406	0.68%	0.059%
16	7.204	693	700	711	rBV	1426449	2164120	32.52%	2.800%
17	7.293	711	715	720	rVB6	15204	21862	0.33%	0.028%
18	7.669	772	779	787	rBV	1445928	2230500	33.51%	2.886%
19	7.746	788	792	797	rVB4	12503	18178	0.27%	0.024%
20	8.387	894	901	914	rBV	1435969	2267400	34.07%	2.934%
21	8.498	917	920	926	rVB	23201	33894	0.51%	0.044%
22	8.834	967	977	988	rBV	1284362	2045735	30.74%	2.647%
23	8.957	993	998	1002	rVV8	6149	11777	0.18%	0.015%
24	9.004	1002	1006	1009	rVB6	5339	8429	0.13%	0.011%
25	9.251	1044	1048	1052	rVB4	11791	16840	0.25%	0.022%
26	9.551	1091	1099	1113	rBV	977638	1589642	23.88%	2.057%
27	10.087	1183	1190	1204	rBV	1437840	2366229	35.55%	3.062%
28	10.275	1217	1222	1228	rBV7	11852	29245	0.44%	0.038%
29	10.463	1247	1254	1264	rBV	1899247	3100509	46.58%	4.012%
30	10.610	1271	1279	1291	rBV	1378029	2343619	35.21%	3.033%
31	10.992	1341	1344	1348	rBV6	5235	7301	0.11%	0.009%
32	11.287	1387	1394	1399	rBV6	6842	14083	0.21%	0.018%
33	11.751	1467	1473	1478	rBV10	6048	9634	0.14%	0.012%
34	11.892	1495	1497	1502	rVB6	5718	8175	0.12%	0.011%
35	12.063	1520	1526	1533	rVB3	30312	51257	0.77%	0.066%
36	12.286	1557	1564	1567	rBV8	4057	9751	0.15%	0.013%

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Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
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37	12.686	1627	1632	1636	rBV7	7802	15247	0.23%	0.020%
38	12.945	1671	1676	1678	rBV4	10427	14712	0.22%	0.019%
39	13.163	1708	1713	1716	rVB2	8650	11404	0.17%	0.015%
40	13.292	1731	1735	1739	rBV6	8175	13503	0.20%	0.017%
41	13.663	1791	1798	1804	rVB10	11993	21078	0.32%	0.027%
42	13.751	1807	1813	1823	rBV	2581757	3535698	53.12%	4.575%
43	14.022	1851	1859	1872	rBV	2904498	4259235	63.99%	5.511%
44	14.245	1892	1897	1901	rVB6	8976	13786	0.21%	0.018%
45	14.333	1905	1912	1919	rVV	2664768	3790070	56.94%	4.904%
46	14.392	1919	1922	1929	rVB9	6410	11099	0.17%	0.014%
47	14.504	1937	1941	1943	rBV5	6090	8111	0.12%	0.010%
48	14.545	1943	1948	1966	rVV	681803	1254020	18.84%	1.623%
49	14.769	1980	1986	1996	rVB	58865	92279	1.39%	0.119%
50	15.151	2044	2051	2053	rBV3	12692	19226	0.29%	0.025%
51	15.333	2076	2082	2090	rBV	3746890	5265166	79.11%	6.813%
52	15.457	2096	2103	2117	rBV	752129	1041113	15.64%	1.347%
53	15.575	2118	2123	2128	rVB	42717	62746	0.94%	0.081%
54	15.710	2144	2146	2153	rVB8	5198	8329	0.13%	0.011%
55	15.910	2176	2180	2186	rVB8	7116	10366	0.16%	0.013%
56	16.027	2194	2200	2201	rBV6	7306	10362	0.16%	0.013%
57	16.075	2205	2208	2213	rVV6	13492	22806	0.34%	0.030%
58	16.127	2213	2217	2225	rVB7	15737	28646	0.43%	0.037%
59	16.251	2233	2238	2240	rBV6	6270	9780	0.15%	0.013%
60	16.592	2290	2296	2298	rBV7	5863	9699	0.15%	0.013%
61	16.651	2301	2306	2310	rVV	33041	46132	0.69%	0.060%
62	16.780	2323	2328	2331	rVB6	8399	11598	0.17%	0.015%
63	16.880	2338	2345	2350	rBV7	14413	31890	0.48%	0.041%
64	16.974	2359	2361	2366	rBV6	5329	8451	0.13%	0.011%
65	17.039	2367	2372	2376	rVV4	16336	22007	0.33%	0.028%
66	17.098	2376	2382	2389	rVV	3082128	4325727	64.99%	5.598%
67	17.198	2393	2399	2410	rVV2	3907124	5558491	83.51%	7.193%
68	17.310	2412	2418	2430	rVB	111858	176715	2.66%	0.229%
69	17.486	2444	2448	2456	rVB8	38624	57654	0.87%	0.075%
70	17.616	2468	2470	2472	rBV3	7426	8119	0.12%	0.011%
71	17.798	2496	2501	2506	rBV	192625	225380	3.39%	0.292%
72	18.010	2533	2537	2542	rBV4	43819	59494	0.89%	0.077%
73	18.080	2547	2549	2553	rVB4	11495	13502	0.20%	0.017%
74	19.033	2707	2711	2715	rVB	54504	71169	1.07%	0.092%
75	19.098	2719	2722	2726	rBV	48957	62927	0.95%	0.081%
76	19.457	2777	2783	2791	rBV	5134704	6655709	100.00%	8.613%

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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

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Title : SVOA CALIBRATION

77	20.951	3033	3037	3045	rBV	377555	605936	9.10%	0.784%
78	21.221	3078	3083	3089	rBV	3818045	4719940	70.92%	6.108%
79	23.421	3450	3457	3464	rBV2	3450169	6485436	97.44%	8.392%
80	23.574	3476	3483	3493	rVB2	2455766	4804050	72.18%	6.216%

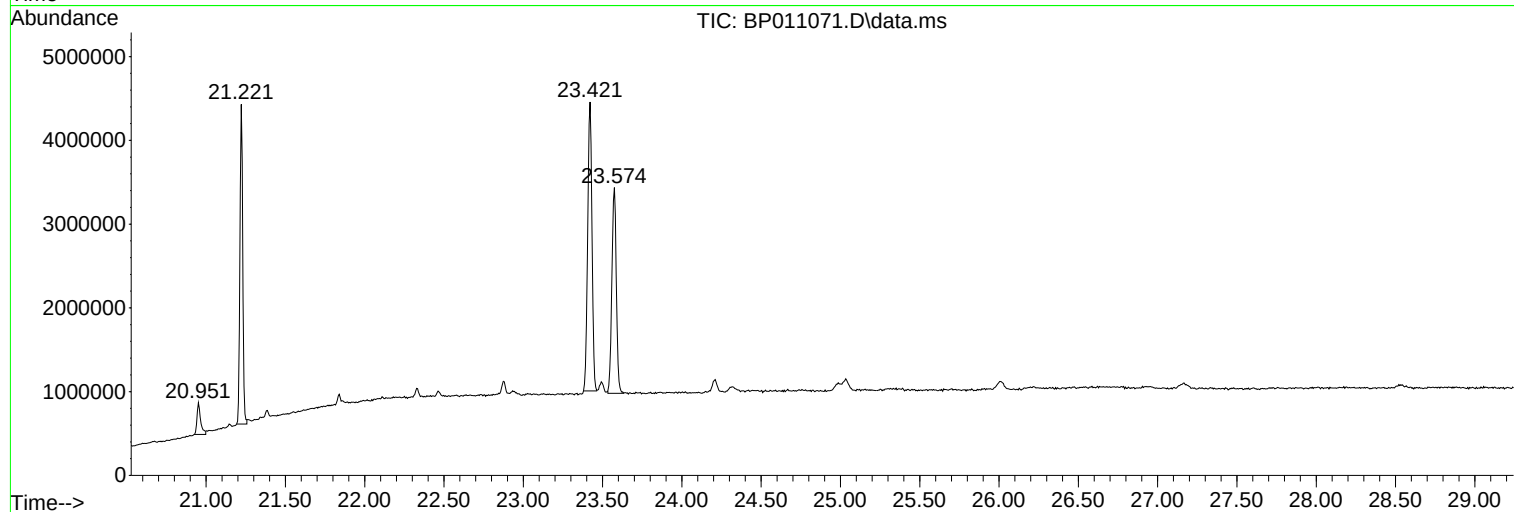
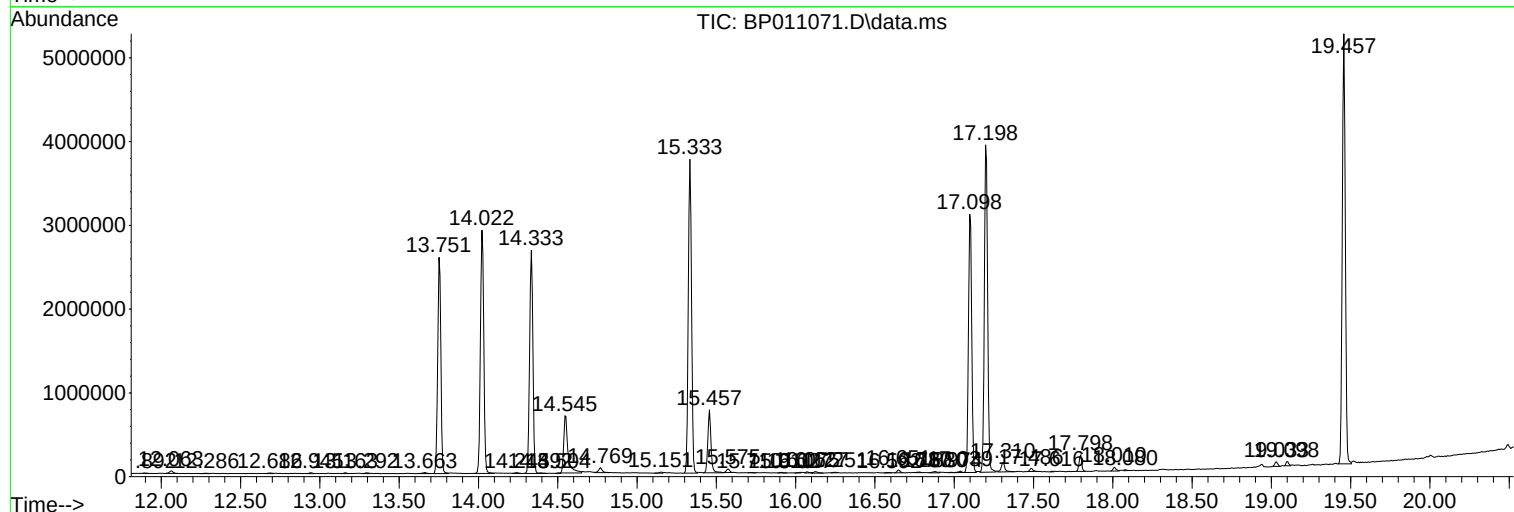
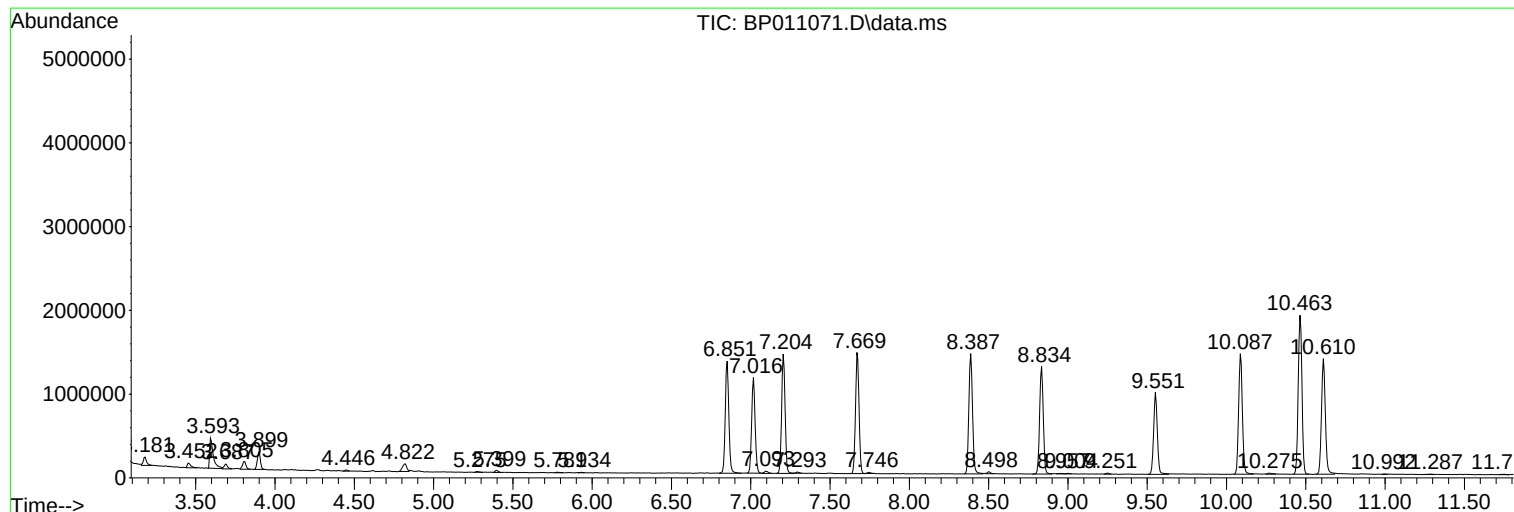
Sum of corrected areas: 77279498

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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



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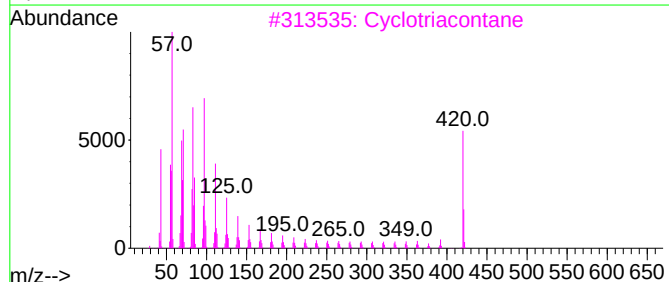
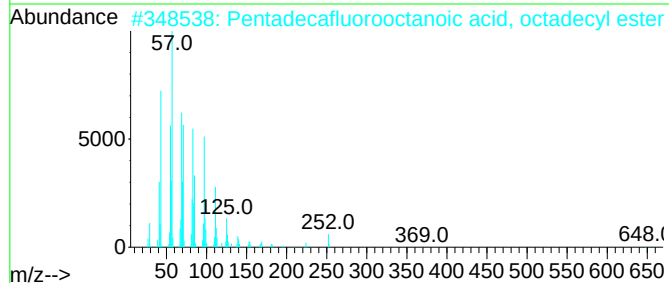
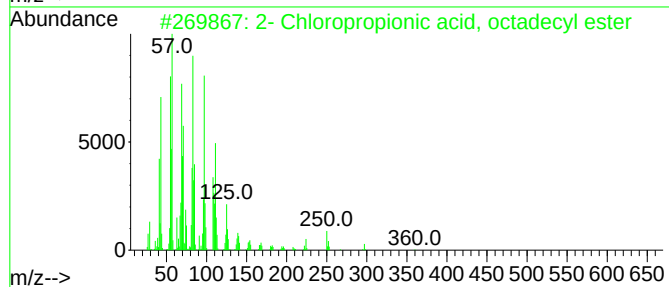
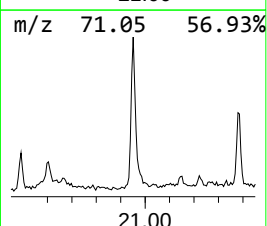
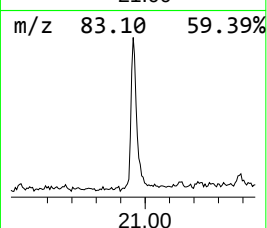
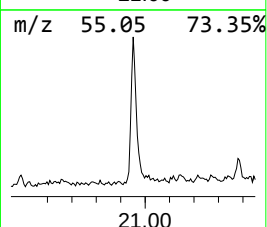
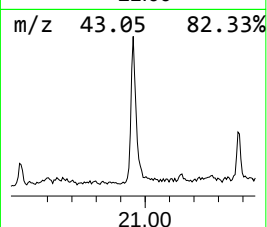
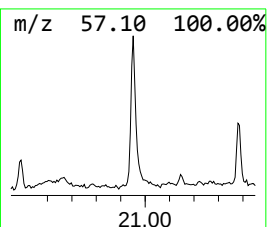
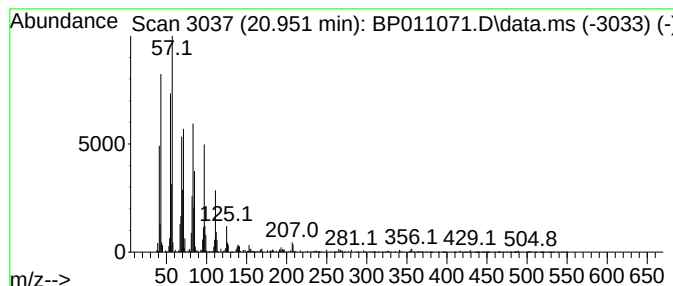
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071422.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 2- Chloropropionic acid, oc... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.951	2.57 ng/ul	605936	Chrysene-d12	21.221

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-	Chloropropionic acid, octadec...	360	C21H41ClO2	088104-31-8	93	
2	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	93		
3	Cyclotriacontane	420	C30H60	000297-35-8	90		
4	Nonadecyl trifluoroacetate	380	C21H39F3O2	1000351-76-3	90		
5	Pentafluoropropionic acid, hepta...	402	C20H35F5O2	959218-78-1	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
2- Chloropropio...	20.951	2.6	ng/u1	605936	5	21.221	4719940	20.0