

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032023\
 Data File : BP014214.D
 Acq On : 22 Mar 2023 11:31
 Operator : CG/JU
 Sample : 01956-02
 Misc :
 ALS Vial : 86 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 E0235

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: BP014214.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.411	34	38	44	rBV	35594	47041	0.69%	0.082%
2	3.534	55	59	63	rBV7	4500	8561	0.13%	0.015%
3	3.852	111	113	131	rBV	162875	293421	4.32%	0.512%
4	4.064	147	149	155	rVB3	7651	11510	0.17%	0.020%
5	4.175	164	168	174	rVB2	12817	19757	0.29%	0.034%
6	7.205	676	683	697	rBV	166184	306352	4.51%	0.534%
7	7.363	704	710	723	rBV	679453	1075199	15.84%	1.875%
8	7.569	739	745	756	rBV	656391	1087174	16.02%	1.896%
9	8.040	818	825	844	rBV	716681	1151411	16.96%	2.008%
10	8.757	939	947	964	rBV	526509	934717	13.77%	1.630%
11	9.216	1018	1025	1040	rBV	898162	1506866	22.20%	2.628%
12	9.599	1084	1090	1093	rBV3	13272	21366	0.31%	0.037%
13	9.940	1140	1148	1163	rBV	731427	1258675	18.54%	2.196%
14	10.487	1233	1241	1260	rBV	940916	1778542	26.20%	3.102%
15	10.863	1297	1305	1314	rBV	1008816	1690444	24.90%	2.949%
16	11.004	1322	1329	1350	rBV	786651	1492490	21.99%	2.603%
17	12.257	1536	1542	1552	rBV5	9994	20829	0.31%	0.036%
18	12.451	1570	1575	1583	rVB2	17931	35333	0.52%	0.062%
19	13.487	1743	1751	1755	rBV6	9077	15776	0.23%	0.028%
20	13.563	1760	1764	1770	rVB2	24409	37415	0.55%	0.065%
21	14.087	1845	1853	1866	rBV	2465888	3400176	50.09%	5.931%
22	14.198	1867	1872	1879	rVB	107655	145134	2.14%	0.253%
23	14.375	1895	1902	1917	rBV	2496040	3689411	54.35%	6.435%
24	14.551	1927	1932	1938	rVB6	11611	17123	0.25%	0.030%
25	14.681	1946	1954	1960	rBV	1688018	2445456	36.03%	4.266%
26	14.739	1960	1964	1986	rVV	354268	990172	14.59%	1.727%
27	14.904	1986	1992	2012	rVB2	114693	356794	5.26%	0.622%
28	15.198	2038	2042	2047	rVB	20038	27267	0.40%	0.048%
29	15.492	2088	2092	2097	rVB2	16568	24429	0.36%	0.043%
30	15.675	2116	2123	2131	rBV	3416264	4882670	71.93%	8.517%
31	15.792	2137	2143	2157	rBV	1336704	1859190	27.39%	3.243%
32	15.916	2161	2164	2169	rVB3	17198	25658	0.38%	0.045%
33	16.039	2179	2185	2191	rBV	74038	112370	1.66%	0.196%
34	16.833	2316	2320	2324	rBV7	4699	8600	0.13%	0.015%
35	17.110	2364	2367	2370	rVB5	7427	10131	0.15%	0.018%
36	17.433	2416	2422	2433	rBV	2127805	3055567	45.02%	5.330%

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37	17.533	2433	2439	2455	rVV	4066985	5867512	86.44%	10.235%
38	18.080	2530	2532	2536	rVB5	8495	8027	0.12%	0.014%
39	18.298	2565	2569	2578	rBV	138109	240997	3.55%	0.420%
40	19.339	2742	2746	2750	rBV2	54409	74938	1.10%	0.131%
41	19.404	2753	2757	2765	rVB	57776	92873	1.37%	0.162%
42	19.522	2774	2777	2781	rBV3	64152	88097	1.30%	0.154%
43	19.757	2811	2817	2823	rBV	5013407	6787644	100.00%	11.840%
44	21.186	3056	3060	3065	rBV	207032	282340	4.16%	0.492%
45	21.510	3110	3115	3123	rBV	2068373	2914846	42.94%	5.084%
46	23.863	3508	3515	3524	rVB2	2444956	4823637	71.06%	8.414%
47	24.027	3537	3543	3551	rVB2	1137809	2305582	33.97%	4.022%

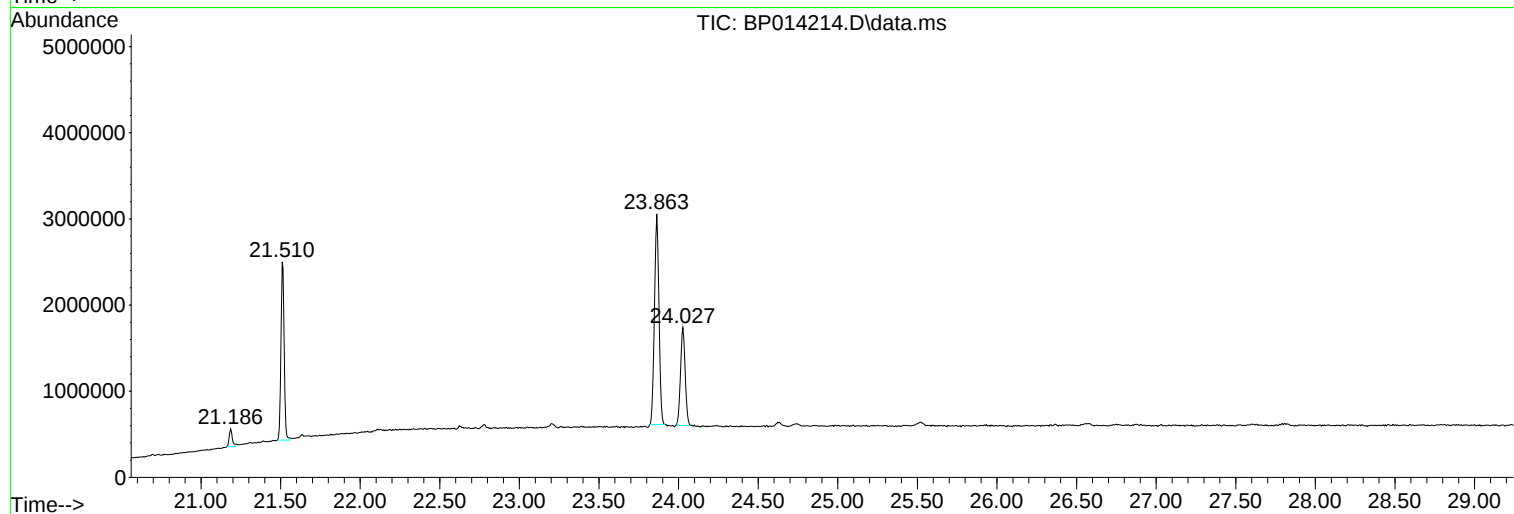
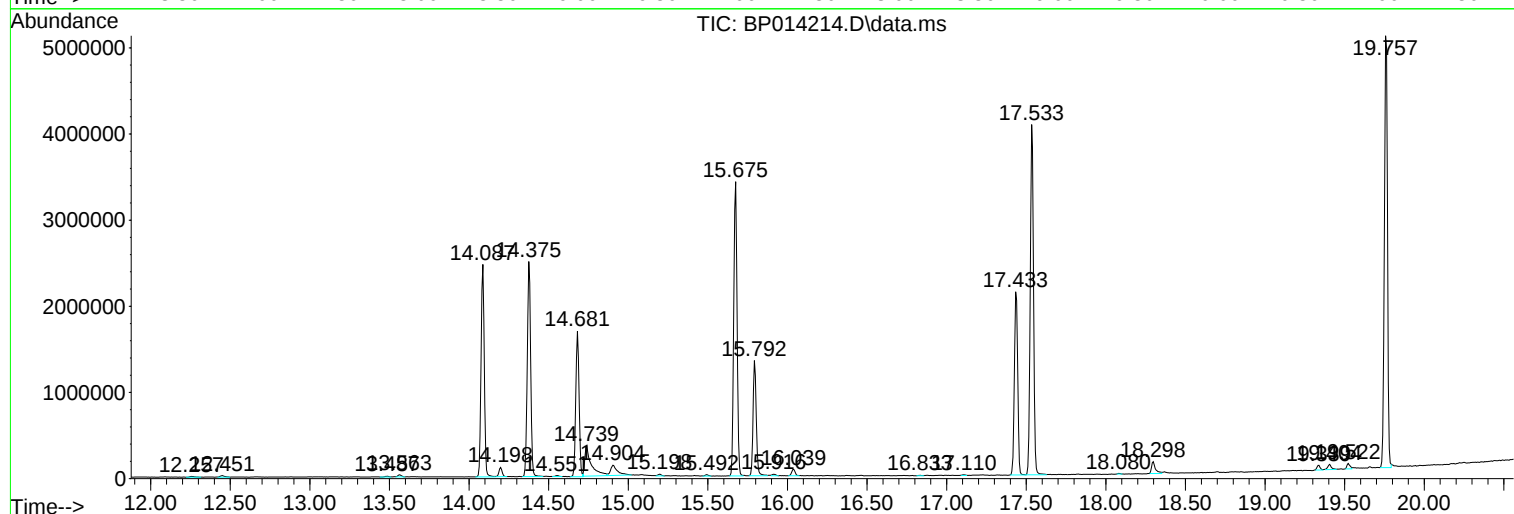
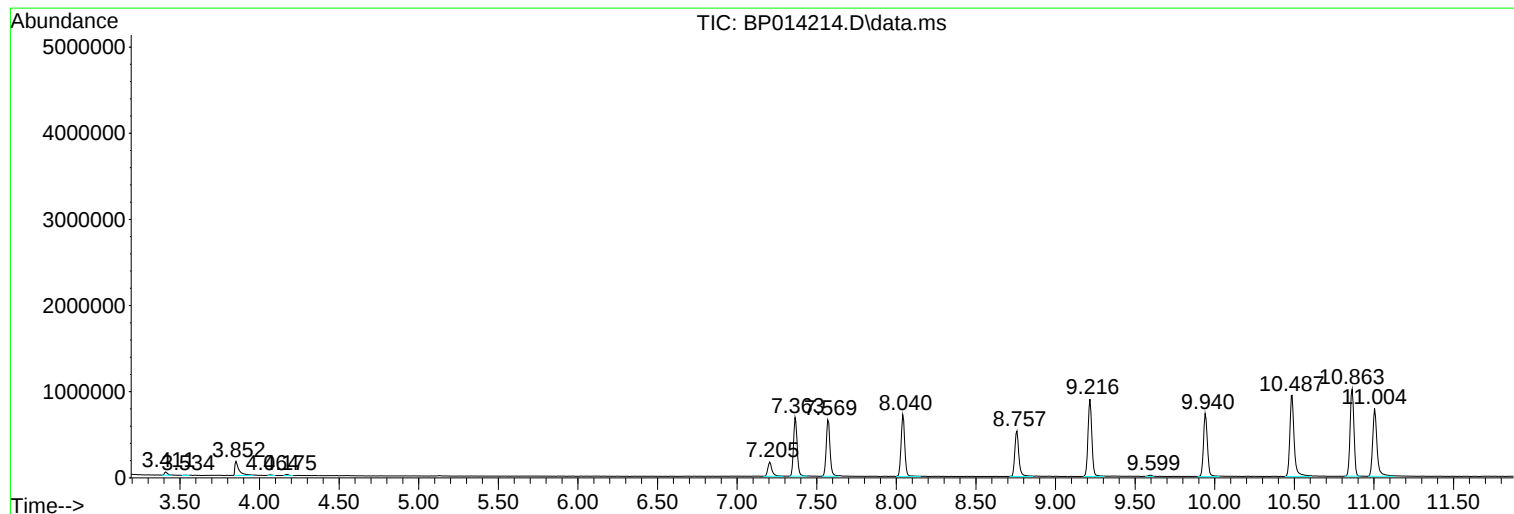
Sum of corrected areas: 57329520

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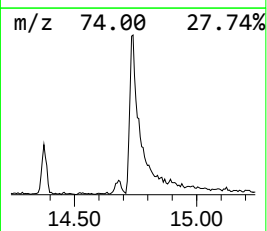
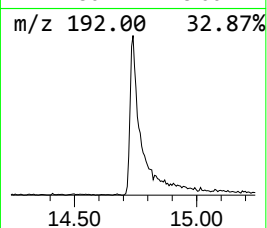
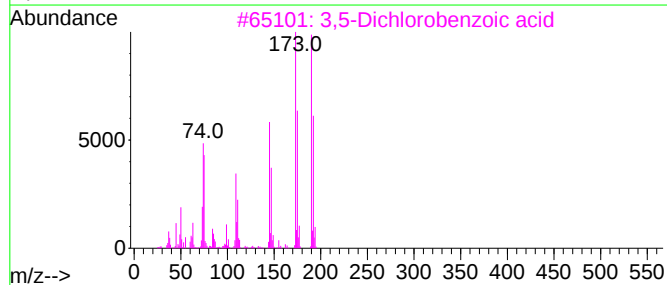
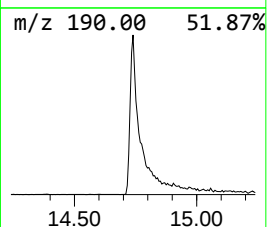
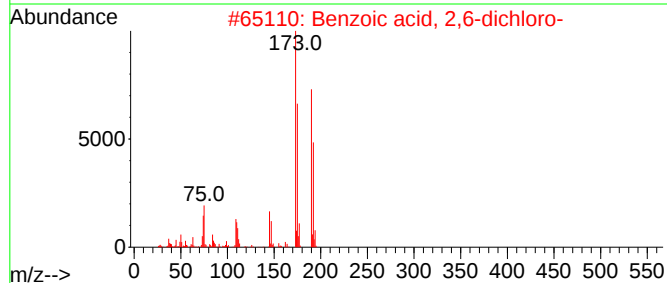
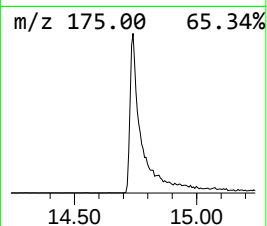
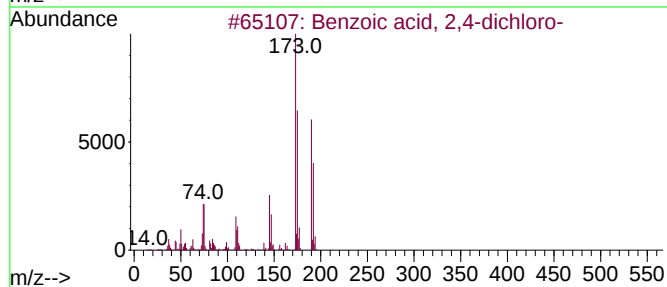
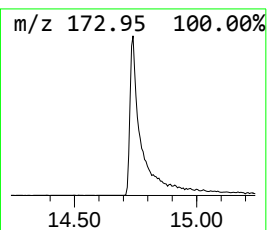
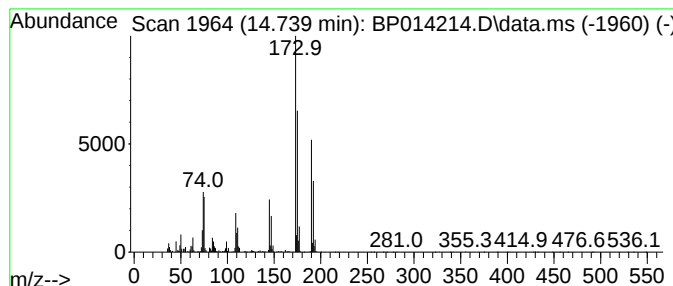
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzoic acid, 2,4-dichloro- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.740	8.10 ng/ul	990172	Acenaphthene-d10	14.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 2,4-dichloro-	190	C7H4Cl2O2	000050-84-0	98
2			Benzoic acid, 2,6-dichloro-	190	C7H4Cl2O2	000050-30-6	97
3			3,5-Dichlorobenzoic acid	190	C7H4Cl2O2	000051-36-5	96
4			Benzoic acid, 3,4-dichloro-	190	C7H4Cl2O2	000051-44-5	96
5			Benzoic acid, 2,4-dichloro-	190	C7H4Cl2O2	000050-84-0	96



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzoic acid, 2...	14.740	8.1	ng/u1	990172	3	14.681	2445460	20.0