

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050123\
 Data File : BP014922.D
 Acq On : 01 May 2023 19:12
 Operator : CG/JU
 Sample : 02528-08DL 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HOA38DL

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-BP042823.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014922.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.434	4	8	11	rBV3	26441	35971	1.26%	0.103%
2	3.464	11	13	16	rVV	27085	35104	1.23%	0.100%
3	3.499	16	19	30	rVB	224514	296203	10.40%	0.846%
4	3.775	61	66	76	rVB8	8116	18416	0.65%	0.053%
5	3.911	87	89	102	rBV	177890	296906	10.43%	0.848%
6	4.152	127	130	136	rVB2	13910	21217	0.75%	0.061%
7	4.252	143	147	153	rVB	23759	32573	1.14%	0.093%
8	4.634	210	212	215	rVB4	3176	3716	0.13%	0.011%
9	4.805	234	241	252	rVB	405165	591293	20.77%	1.689%
10	5.205	306	309	316	rVB4	6883	11618	0.41%	0.033%
11	5.699	390	393	398	rVB2	6692	8443	0.30%	0.024%
12	5.793	404	409	411	rBV6	3198	4914	0.17%	0.014%
13	5.822	411	414	421	rVB	8625	14210	0.50%	0.041%
14	6.293	490	494	497	rVB6	2207	3083	0.11%	0.009%
15	6.546	532	537	544	rVB2	10831	20121	0.71%	0.057%
16	7.311	657	667	680	rBV	205035	353359	12.41%	1.010%
17	7.487	690	697	708	rBV	370862	597394	20.98%	1.707%
18	7.693	725	732	743	rBV	400241	641768	22.54%	1.833%
19	7.975	775	780	786	rBV8	2379	4150	0.15%	0.012%
20	8.169	804	813	820	rBV	780747	1251256	43.95%	3.575%
21	8.222	820	822	828	rVB3	11771	18049	0.63%	0.052%
22	8.534	870	875	878	rBV7	3618	6599	0.23%	0.019%
23	8.581	882	883	890	rVB7	1890	2904	0.10%	0.008%
24	8.875	926	933	943	rBV	334696	561412	19.72%	1.604%
25	9.346	1004	1013	1026	rBV	367907	637818	22.40%	1.822%
26	9.428	1026	1027	1033	rVB5	5443	7726	0.27%	0.022%
27	9.681	1066	1070	1075	rBV8	2120	3986	0.14%	0.011%
28	9.752	1076	1082	1090	rVB2	27598	45261	1.59%	0.129%
29	10.075	1130	1137	1150	rBV	270206	467034	16.41%	1.334%
30	10.210	1159	1160	1168	rVB8	2549	3757	0.13%	0.011%
31	10.369	1182	1187	1192	rBV9	1820	3689	0.13%	0.011%
32	10.616	1221	1229	1241	rBV	483566	823382	28.92%	2.352%
33	10.722	1245	1247	1251	rVB4	2001	2877	0.10%	0.008%
34	11.005	1286	1295	1306	rBV	1057444	1801946	63.30%	5.148%
35	11.140	1311	1318	1334	rVV2	159065	319302	11.22%	0.912%
36	11.393	1359	1361	1365	rVB5	2803	3036	0.11%	0.009%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
 Title : SVOA CALIBRATION

37	11.781	1420	1427	1428	rBV6	4221	5347	0.19%	0.015%
38	11.816	1428	1433	1439	rVB9	5597	11681	0.41%	0.033%
39	12.134	1483	1487	1488	rBV4	2559	3206	0.11%	0.009%
40	12.257	1502	1508	1513	rVB5	6186	10274	0.36%	0.029%
41	12.310	1516	1517	1527	rVB10	3220	5025	0.18%	0.014%
42	12.587	1559	1564	1569	rVB4	8168	14313	0.50%	0.041%
43	12.793	1597	1599	1603	rVB5	2948	3643	0.13%	0.010%
44	13.175	1659	1664	1667	rBV7	2152	3970	0.14%	0.011%
45	13.634	1737	1742	1743	rBV4	4970	6462	0.23%	0.018%
46	13.704	1749	1754	1757	rBV	17704	27688	0.97%	0.079%
47	13.787	1766	1768	1772	rVB5	2880	3626	0.13%	0.010%
48	13.904	1786	1788	1792	rVB4	2424	3113	0.11%	0.009%
49	13.946	1792	1795	1797	rBV4	2507	3286	0.12%	0.009%
50	13.975	1797	1800	1804	rVV6	2891	5413	0.19%	0.015%
51	14.016	1804	1807	1811	rVB6	1659	3235	0.11%	0.009%
52	14.216	1834	1841	1846	rBV	1033163	1421059	49.92%	4.060%
53	14.287	1846	1853	1857	rBV3	59227	181206	6.37%	0.518%
54	14.510	1885	1891	1912	rVB	1197073	1792005	62.95%	5.120%
55	14.698	1921	1923	1930	rVB7	4919	7742	0.27%	0.022%
56	14.816	1936	1943	1950	rBV	1541931	2310774	81.17%	6.602%
57	14.898	1955	1957	1962	rVB6	4830	4973	0.17%	0.014%
58	14.993	1966	1973	1991	rBV	98037	207812	7.30%	0.594%
59	15.216	2005	2011	2016	rBV10	5784	12692	0.45%	0.036%
60	15.257	2016	2018	2023	rVB6	4257	5994	0.21%	0.017%
61	15.604	2068	2077	2078	rBV9	3775	7171	0.25%	0.020%
62	15.645	2082	2084	2088	rVB5	3257	3351	0.12%	0.010%
63	15.710	2094	2095	2102	rVB7	2431	2874	0.10%	0.008%
64	15.804	2102	2111	2124	rBV	1476733	2175351	76.41%	6.215%
65	15.916	2124	2130	2140	rVV	345241	476599	16.74%	1.362%
66	16.045	2148	2152	2157	rVB7	8000	12898	0.45%	0.037%
67	16.169	2167	2173	2181	rBV	693803	941760	33.08%	2.691%
68	16.640	2248	2253	2255	rBV5	2650	5021	0.18%	0.014%
69	16.734	2265	2269	2275	rBV	96508	139845	4.91%	0.400%
70	16.869	2288	2292	2293	rBV3	3223	3429	0.12%	0.010%
71	16.945	2302	2305	2307	rBV3	4150	4277	0.15%	0.012%
72	17.098	2328	2331	2332	rBV3	3933	3166	0.11%	0.009%
73	17.245	2353	2356	2359	rBV5	3552	4425	0.16%	0.013%
74	17.322	2364	2369	2371	rBV6	4758	7773	0.27%	0.022%
75	17.357	2371	2375	2380	rVV8	6973	12813	0.45%	0.037%
76	17.469	2391	2394	2395	rBV3	2622	3149	0.11%	0.009%

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Title : SVOA CALIBRATION

77	17.569	2405	2411	2419	rBV	1797415	2622397	92.11%	7.492%
78	17.669	2422	2428	2441	rVV	1591324	2213796	77.76%	6.325%
79	18.428	2552	2557	2565	rBV	125756	184519	6.48%	0.527%
80	18.510	2567	2571	2579	rVB5	37430	64065	2.25%	0.183%
81	19.533	2742	2745	2748	rBV	21728	26583	0.93%	0.076%
82	19.651	2762	2765	2770	rBV3	63295	72253	2.54%	0.206%
83	19.886	2800	2805	2818	rBV	1961030	2718358	95.49%	7.766%
84	21.327	3047	3050	3056	rBV3	103735	147397	5.18%	0.421%
85	21.651	3100	3105	3115	rBV2	1918863	2731329	95.94%	7.803%
86	24.086	3513	3519	3535	rVB2	1149701	2566256	90.14%	7.332%
87	24.263	3542	3549	3563	rVB2	1275806	2846875	100.00%	8.133%

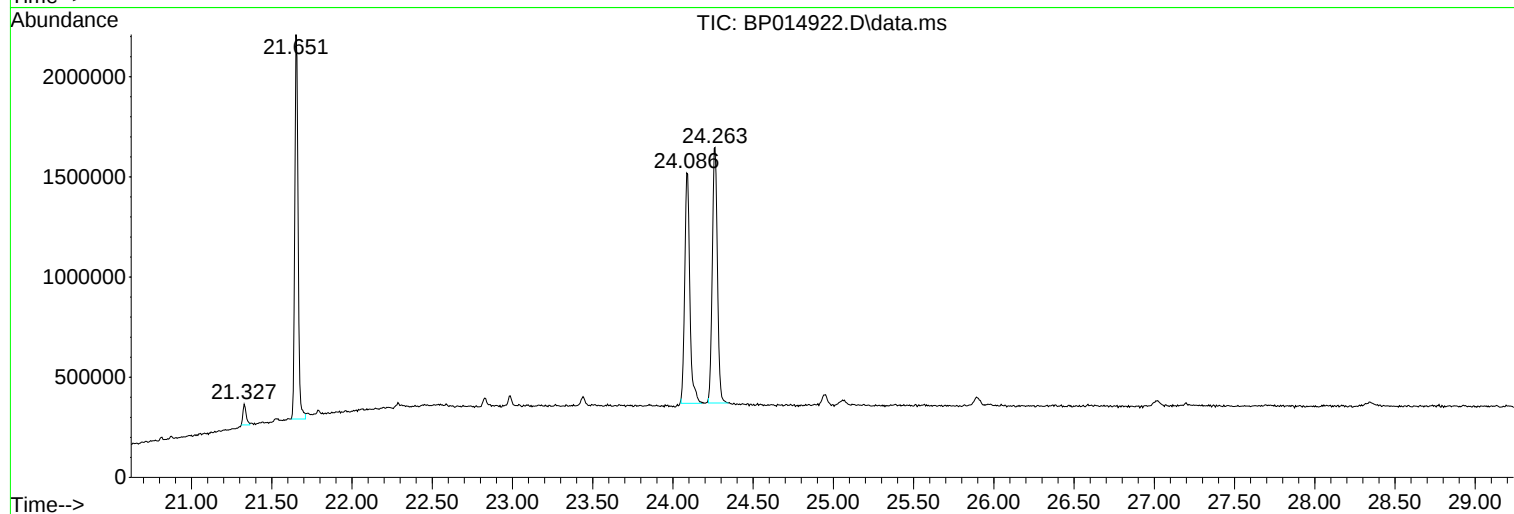
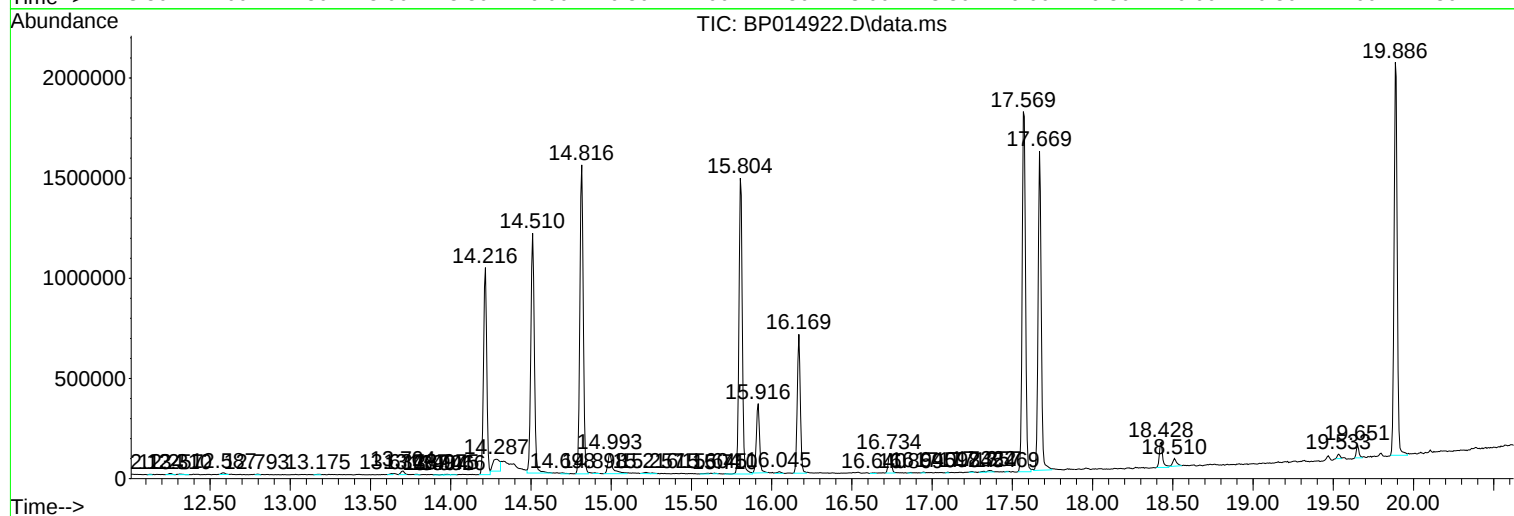
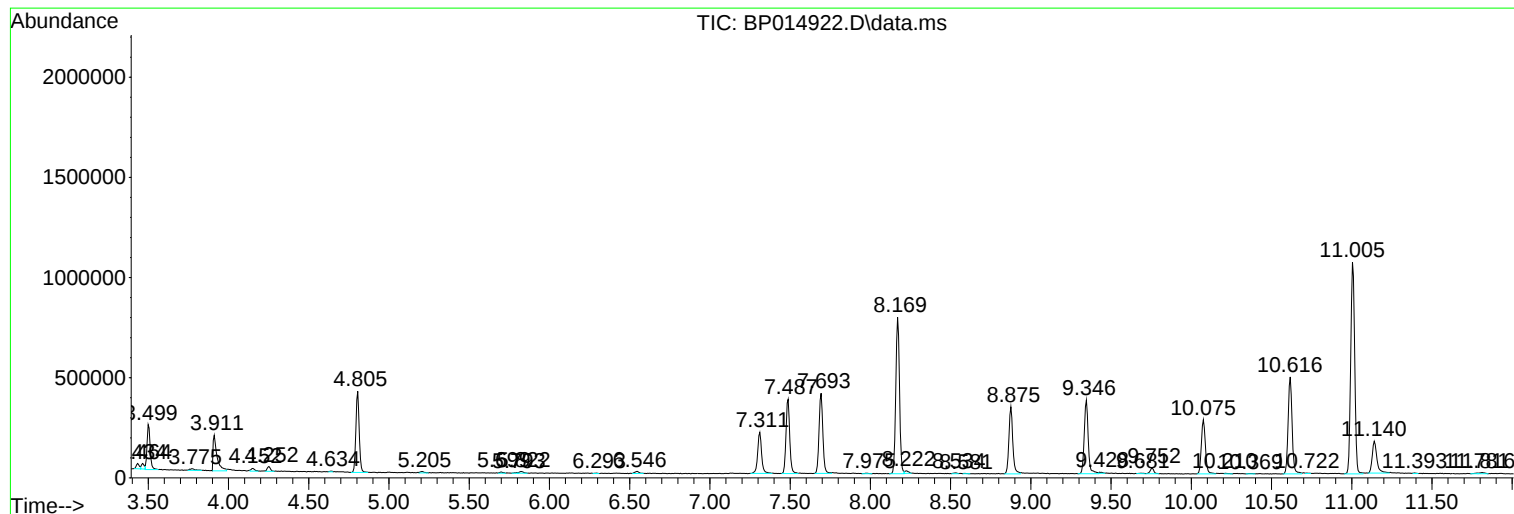
Sum of corrected areas: 35002762

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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



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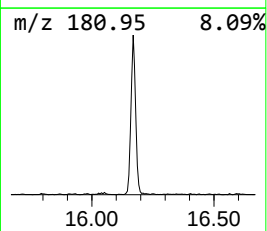
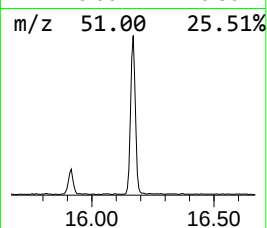
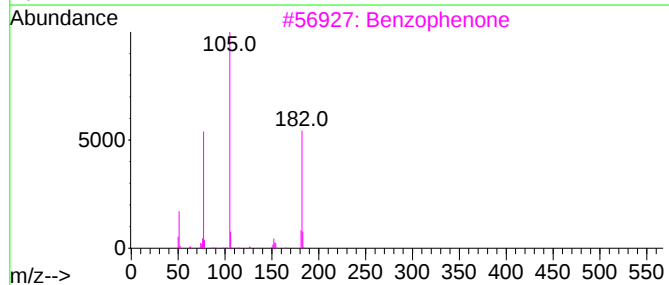
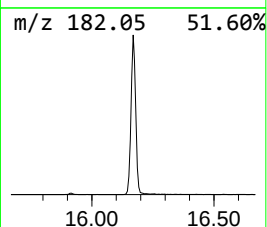
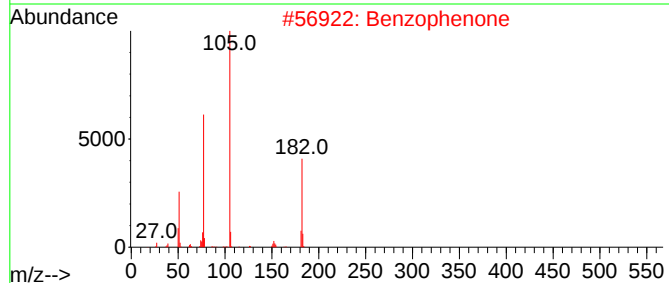
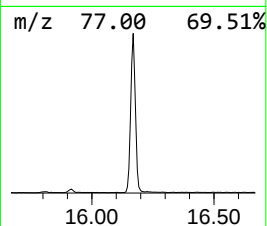
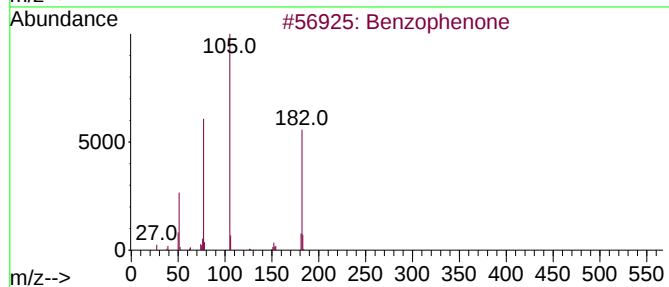
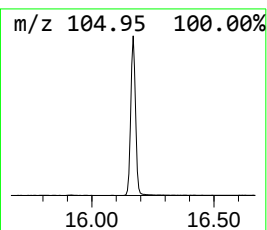
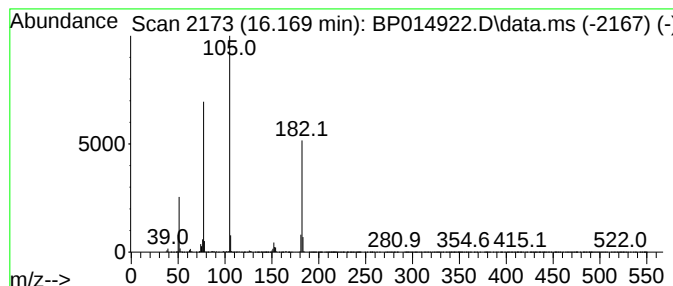
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042823.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.169	8.15 ng/ul	941760	Acenaphthene-d10	14.816

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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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
Benzophenone	16.169	8.2	ng/u1	941760	3	14.816	2310770	20.0