

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040523\
 Data File : BP014577.D
 Acq On : 06 Apr 2023 03:15
 Operator : CG/JU
 Sample : 02194-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHDW0

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014577.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	20	24	28	rBV	120297	161250	3.56%	0.399%
2	3.364	28	30	34	rVV	35971	45779	1.01%	0.113%
3	3.405	34	37	42	rVB2	18207	23864	0.53%	0.059%
4	3.822	105	108	120	rBV	48719	116709	2.58%	0.289%
5	4.016	136	141	149	rVB2	17202	26091	0.58%	0.065%
6	4.122	155	159	164	rVB2	10581	16922	0.37%	0.042%
7	4.228	174	177	181	rBV6	8468	11697	0.26%	0.029%
8	4.652	243	249	264	rVB	902002	1340629	29.58%	3.314%
9	4.852	280	283	287	rVB6	3030	5019	0.11%	0.012%
10	5.052	313	317	323	rVB2	10127	17058	0.38%	0.042%
11	6.369	536	541	548	rVB5	6896	10908	0.24%	0.027%
12	7.081	657	662	664	rBV5	3974	5013	0.11%	0.012%
13	7.163	670	676	687	rBV	80177	193418	4.27%	0.478%
14	7.310	695	701	715	rBV	475744	778981	17.19%	1.926%
15	7.516	729	736	753	rBV	413234	765522	16.89%	1.893%
16	7.981	807	815	826	rBV	498569	860234	18.98%	2.127%
17	8.710	932	939	959	rBV2	285479	624899	13.79%	1.545%
18	9.163	1009	1016	1037	rBV	586185	1069082	23.59%	2.643%
19	9.528	1074	1078	1083	rVB4	7792	11978	0.26%	0.030%
20	9.887	1131	1139	1156	rBV	462455	859694	18.97%	2.125%
21	10.434	1225	1232	1258	rBV	493556	1135273	25.05%	2.807%
22	10.804	1288	1295	1310	rBV	733482	1273229	28.10%	3.148%
23	10.957	1314	1321	1343	rBV	297527	774740	17.10%	1.915%
24	12.416	1563	1569	1580	rBV8	8928	20365	0.45%	0.050%
25	13.434	1739	1742	1747	rVB6	3763	6113	0.13%	0.015%
26	13.522	1753	1757	1764	rVB7	6362	13243	0.29%	0.033%
27	14.045	1838	1846	1862	rBV	1626474	2362523	52.13%	5.841%
28	14.151	1862	1864	1868	rVB5	6015	8633	0.19%	0.021%
29	14.334	1888	1895	1913	rBV	1686115	2515340	55.51%	6.218%
30	14.639	1939	1947	1959	rBV	1266530	1881870	41.53%	4.652%
31	14.869	1982	1986	1990	rBV6	6108	9518	0.21%	0.024%
32	14.916	1990	1994	1995	rBV2	16519	19985	0.44%	0.049%
33	15.416	2076	2079	2083	rBV2	3914	5832	0.13%	0.014%
34	15.457	2083	2086	2092	rVB4	6952	8390	0.19%	0.021%
35	15.634	2109	2116	2132	rBV	2200096	3288848	72.58%	8.131%
36	15.763	2132	2138	2151	rBV	728213	1137184	25.09%	2.811%

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 Title : SVOA CALIBRATION

37	16.004	2173	2179	2190	rBV	194996	310766	6.86%	0.768%
38	16.328	2232	2234	2240	rVB7	3843	6338	0.14%	0.016%
39	16.563	2270	2274	2279	rBV4	6193	10918	0.24%	0.027%
40	17.398	2410	2416	2426	rBV	1652356	2297995	50.71%	5.681%
41	17.498	2427	2433	2449	rVB	2623946	3867989	85.36%	9.562%
42	18.275	2561	2565	2572	rBV6	23918	49583	1.09%	0.123%
43	19.304	2737	2740	2746	rBV2	37675	48288	1.07%	0.119%
44	19.374	2748	2752	2760	rVV	35028	61206	1.35%	0.151%
45	19.727	2806	2812	2828	rBV	3360101	4531635	100.00%	11.203%
46	21.486	3106	3111	3125	rBV	1415930	2268942	50.07%	5.609%
47	23.821	3501	3508	3521	rVB2	1735405	3644125	80.42%	9.009%
48	23.986	3530	3536	3548	rVB2	912244	1945952	42.94%	4.811%

Sum of corrected areas: 40449570

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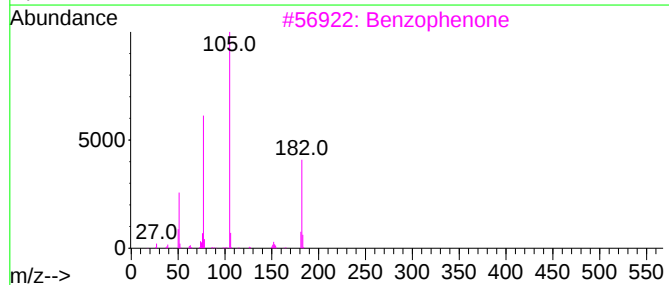
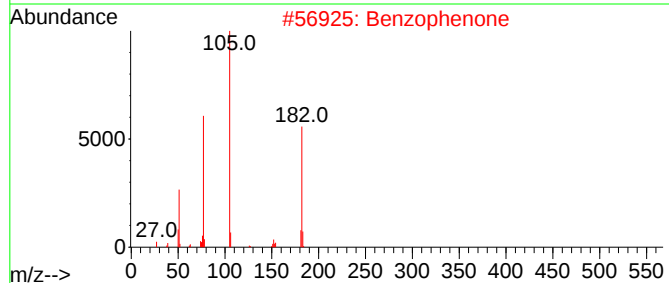
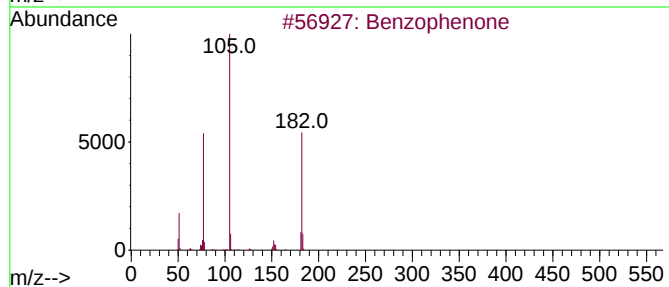
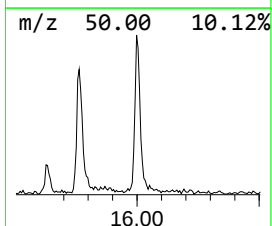
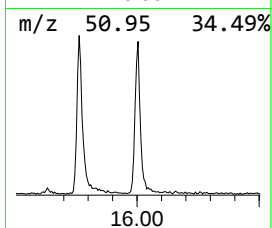
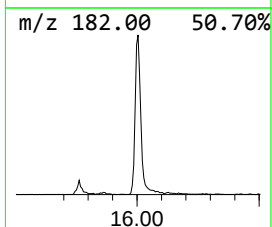
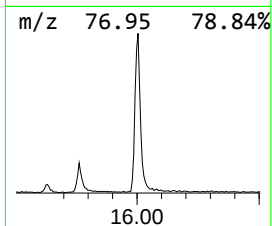
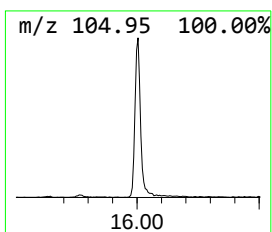
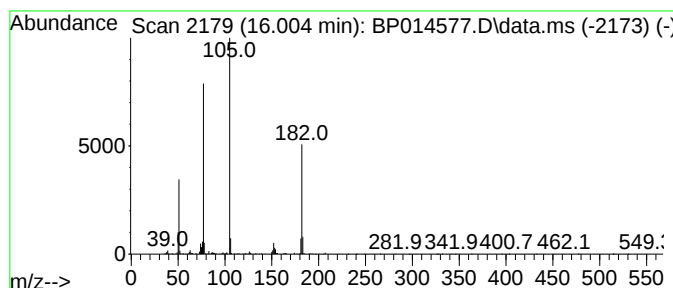
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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.
16.004	3.30 ng/ul	310766	Acenaphthene-d10	14.639

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	93
5			Benzophenone	182	C13H10O	000119-61-9	91



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.004	3.3	ng/u1	310766	3	14.639	1881870	20.0