

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101424\
 Data File : BP022397.D
 Acq On : 15 Oct 2024 20:21
 Operator : RC/JU
 Sample : P4349-16
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EOAG0

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022397.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.223	36	40	45	rVB	15767	20479	0.60%	0.076%
2	3.640	107	111	123	rBV	70744	114996	3.40%	0.427%
3	3.952	161	164	169	rVB2	5354	7839	0.23%	0.029%
4	6.881	652	662	676	rVB	92269	160175	4.73%	0.595%
5	7.034	681	688	697	rBV	242822	390893	11.55%	1.453%
6	7.234	713	722	735	rBV	306733	510719	15.09%	1.898%
7	7.693	792	800	809	rBV	300346	476407	14.07%	1.770%
8	8.393	913	919	928	rBV	267247	444945	13.14%	1.653%
9	8.834	986	994	1009	rBV	323207	589302	17.41%	2.190%
10	9.246	1059	1064	1071	rVB5	4693	8320	0.25%	0.031%
11	9.393	1081	1089	1096	rVB4	7407	12635	0.37%	0.047%
12	9.546	1106	1115	1126	rBV	298047	526128	15.54%	1.955%
13	10.081	1198	1206	1220	rBV	467634	831695	24.57%	3.091%
14	10.452	1260	1269	1282	rBV	390961	683298	20.19%	2.539%
15	10.587	1284	1292	1306	rBV	287408	570574	16.86%	2.120%
16	12.040	1533	1539	1549	rBV2	8930	17333	0.51%	0.064%
17	13.710	1816	1823	1836	rBV	1005980	1333579	39.40%	4.956%
18	13.998	1863	1872	1888	rBV	981867	1542673	45.57%	5.732%
19	14.310	1917	1925	1933	rBV	674676	1073286	31.71%	3.988%
20	14.534	1956	1963	1977	rBV	83373	175253	5.18%	0.651%
21	14.687	1985	1989	1992	rBV6	2217	3901	0.12%	0.014%
22	14.746	1995	1999	2007	rVV10	4237	8684	0.26%	0.032%
23	15.316	2088	2096	2110	rBV	1155609	2066350	61.04%	7.678%
24	15.457	2113	2120	2132	rVV	521182	703311	20.78%	2.613%
25	15.557	2134	2137	2147	rVB4	7294	13294	0.39%	0.049%
26	15.693	2152	2160	2170	rBV	106460	186119	5.50%	0.692%
27	16.045	2212	2220	2227	rBV	24485	42881	1.27%	0.159%
28	16.704	2327	2332	2338	rBV9	2824	6839	0.20%	0.025%
29	16.787	2341	2346	2352	rBV10	2882	6539	0.19%	0.024%
30	17.116	2395	2402	2409	rBV	970982	1479512	43.71%	5.498%
31	17.216	2412	2419	2429	rVV	1959572	2460590	72.69%	9.143%
32	17.422	2449	2454	2457	rBV7	3285	6002	0.18%	0.022%
33	18.051	2556	2561	2568	rBV2	36253	58703	1.73%	0.218%
34	18.551	2644	2646	2652	rVB6	4807	5718	0.17%	0.021%
35	19.139	2741	2746	2751	rBV2	21397	32853	0.97%	0.122%
36	19.216	2754	2759	2763	rBV	31393	46384	1.37%	0.172%

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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

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Title : SVOA CALIBRATION

37	19.375	2783	2786	2793	rVB6	18921	28299	0.84%	0.105%
38	19.539	2811	2814	2817	rBV5	8563	8699	0.26%	0.032%
39	19.598	2817	2824	2831	rBV	2348669	3100545	91.59%	11.521%
40	21.210	3095	3098	3103	rVB3	33003	49881	1.47%	0.185%
41	21.539	3148	3154	3163	rVB	1139152	1792950	52.97%	6.663%
42	24.610	3666	3676	3689	rVB	1126926	3385127	100.00%	12.579%
43	24.827	3705	3713	3726	rVB	693846	1927357	56.94%	7.162%

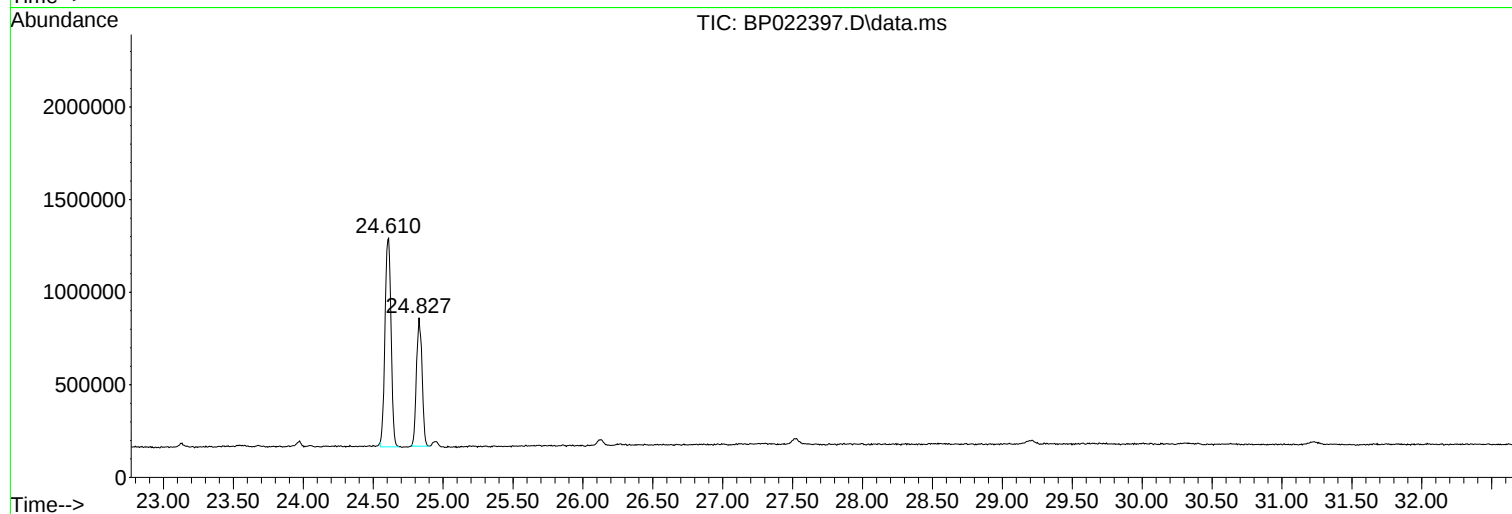
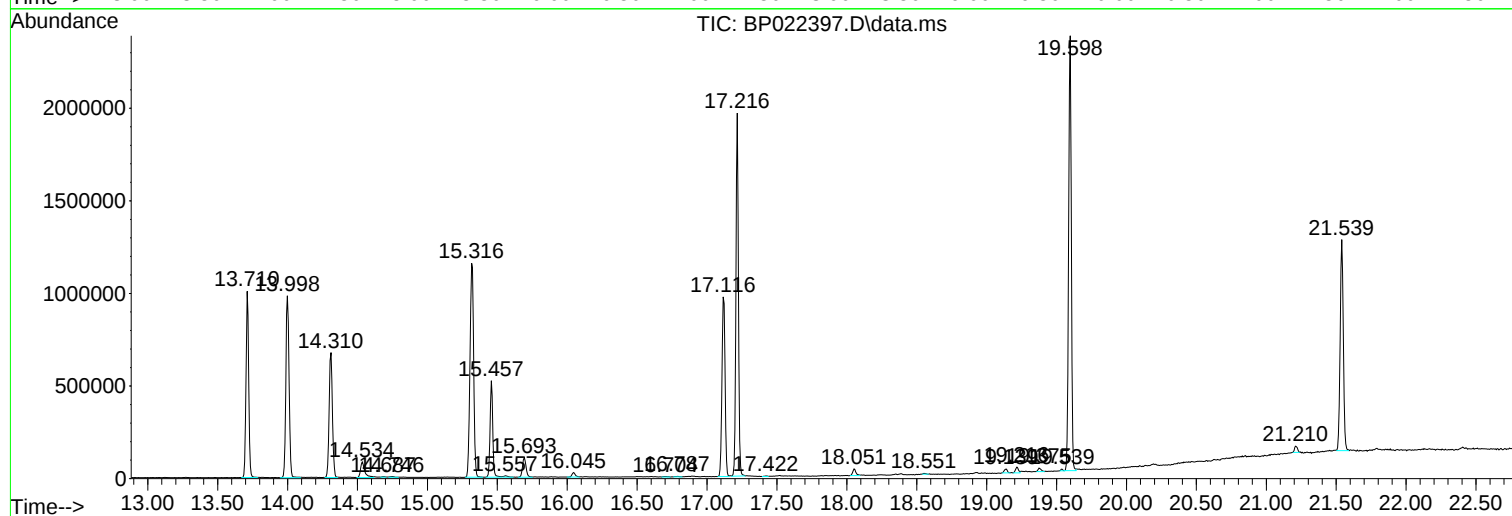
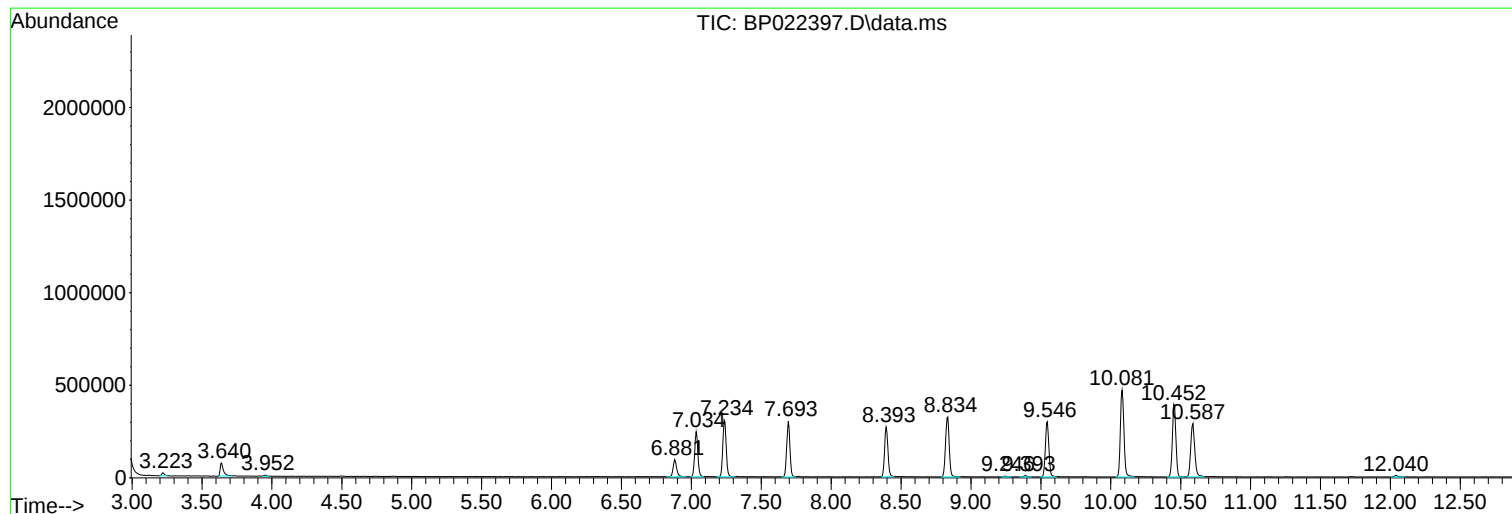
Sum of corrected areas: 26911067

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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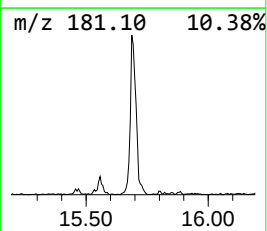
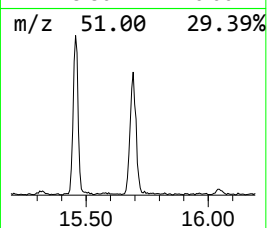
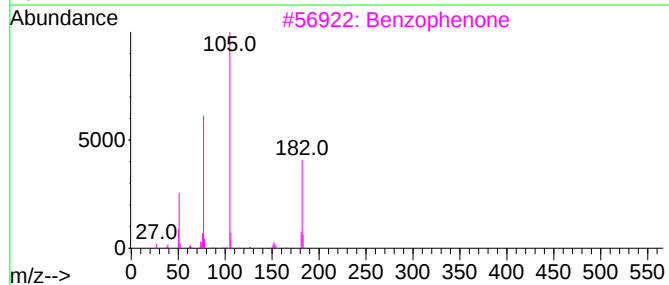
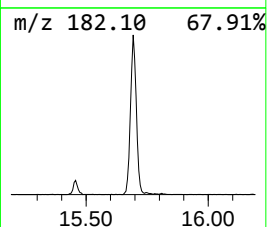
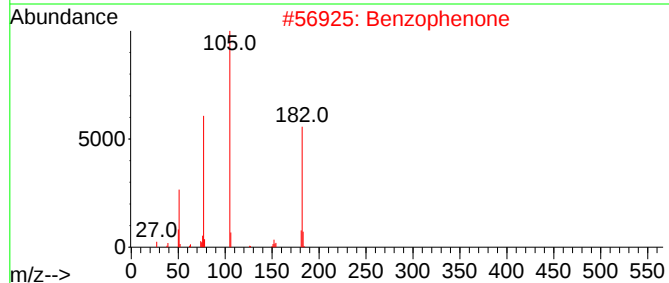
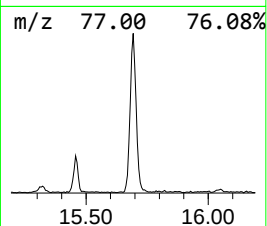
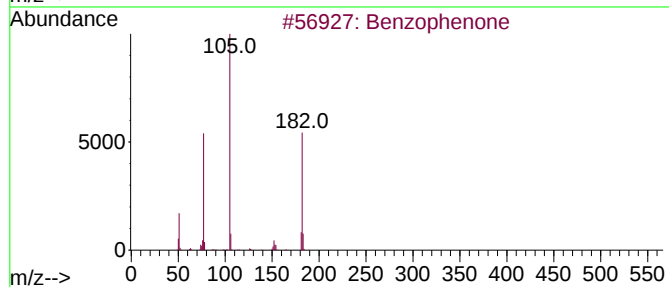
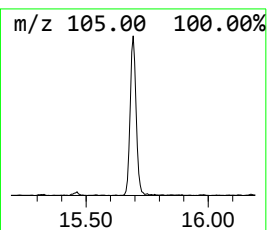
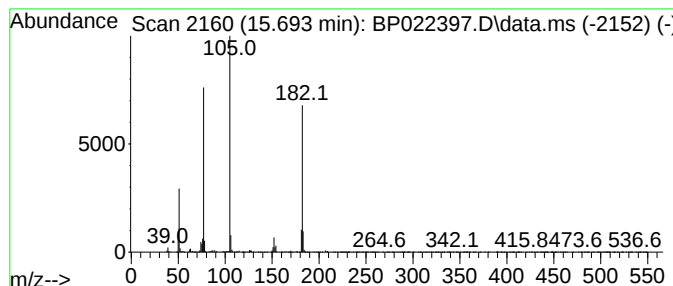
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.693	3.47 ng/ul	186119	Acenaphthene-d10	14.310

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	91
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	15.693	3.5	ng/u1	186119	3	14.310	1073290	20.0