

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
 Data File : BP022813.D
 Acq On : 01 Nov 2024 20:52
 Operator : RC/JU
 Sample : P4627-04
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0CR2

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: BP022813.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.175	29	32	38	rVB	12849	16601	0.56%	0.081%
2	3.593	100	103	118	rBV	52560	99375	3.33%	0.486%
3	4.393	235	239	248	rVB	73464	98632	3.31%	0.482%
4	5.322	392	397	405	rBV5	4322	9802	0.33%	0.048%
5	5.681	453	458	463	rBV7	2454	3661	0.12%	0.018%
6	6.311	557	565	570	rBV6	5273	8629	0.29%	0.042%
7	6.828	648	653	669	rBV	72183	142944	4.79%	0.699%
8	6.969	670	677	689	rVV	205301	302861	10.15%	1.481%
9	7.175	704	712	723	rBV	225806	406419	13.63%	1.988%
10	7.269	725	728	736	rVB7	2879	5757	0.19%	0.028%
11	7.628	783	789	798	rBV	209458	337608	11.32%	1.651%
12	7.693	798	800	808	rVB9	2366	5083	0.17%	0.025%
13	8.340	902	910	924	rBV2	235895	408790	13.71%	1.999%
14	8.775	976	984	993	rBV	275491	464052	15.56%	2.269%
15	9.169	1047	1051	1057	rVB7	3994	6307	0.21%	0.031%
16	9.487	1097	1105	1119	rBV2	226243	419514	14.07%	2.052%
17	10.028	1189	1197	1216	rBV	303447	651016	21.83%	3.184%
18	10.387	1250	1258	1270	rBV	298819	458013	15.36%	2.240%
19	10.534	1276	1283	1294	rBV	254072	470649	15.78%	2.302%
20	11.093	1373	1378	1384	rVB9	1627	3030	0.10%	0.015%
21	11.657	1468	1474	1479	rBV6	2763	5140	0.17%	0.025%
22	12.004	1526	1533	1537	rBV5	3997	8262	0.28%	0.040%
23	12.851	1672	1677	1682	rVB9	1302	3054	0.10%	0.015%
24	13.663	1807	1815	1825	rBV	642062	1051800	35.27%	5.144%
25	13.945	1857	1863	1871	rBV	842741	1114938	37.38%	5.453%
26	14.257	1910	1916	1928	rVB	509028	682095	22.87%	3.336%
27	14.534	1956	1963	1977	rBV4	41714	132560	4.44%	0.648%
28	14.722	1993	1995	2004	rVB10	3259	7212	0.24%	0.035%
29	15.275	2081	2089	2096	rBV	1131149	1566718	52.53%	7.662%
30	15.322	2096	2097	2104	rVB6	4850	7172	0.24%	0.035%
31	15.422	2107	2114	2125	rBV	268915	476506	15.98%	2.330%
32	15.639	2145	2151	2161	rVB	61632	85810	2.88%	0.420%
33	16.734	2335	2337	2345	rVB8	2388	3908	0.13%	0.019%
34	17.063	2387	2393	2402	rBV	644253	899133	30.15%	4.397%
35	17.163	2403	2410	2420	rBV	1367047	1940929	65.08%	9.492%
36	17.363	2442	2444	2449	rVB4	2876	3701	0.12%	0.018%

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Title : SVOA CALIBRATION

37	17.751	2507	2510	2515	rBV6	6663	9685	0.32%	0.047%
38	17.998	2547	2552	2558	rBV2	37307	55289	1.85%	0.270%
39	18.075	2561	2565	2569	rVB	5221	7747	0.26%	0.038%
40	18.310	2603	2605	2609	rVB4	3296	3339	0.11%	0.016%
41	19.104	2735	2740	2745	rBV2	17226	27731	0.93%	0.136%
42	19.180	2748	2753	2756	rBV	19360	26676	0.89%	0.130%
43	19.333	2775	2779	2785	rBV5	26262	38795	1.30%	0.190%
44	19.557	2811	2817	2826	rBV	1930331	2542206	85.24%	12.433%
45	21.504	3142	3148	3159	rVB	645491	1178264	39.51%	5.762%
46	24.545	3655	3665	3681	rVB	1157087	2982516	100.00%	14.586%
47	24.763	3693	3702	3714	rVB	451086	1268006	42.51%	6.201%

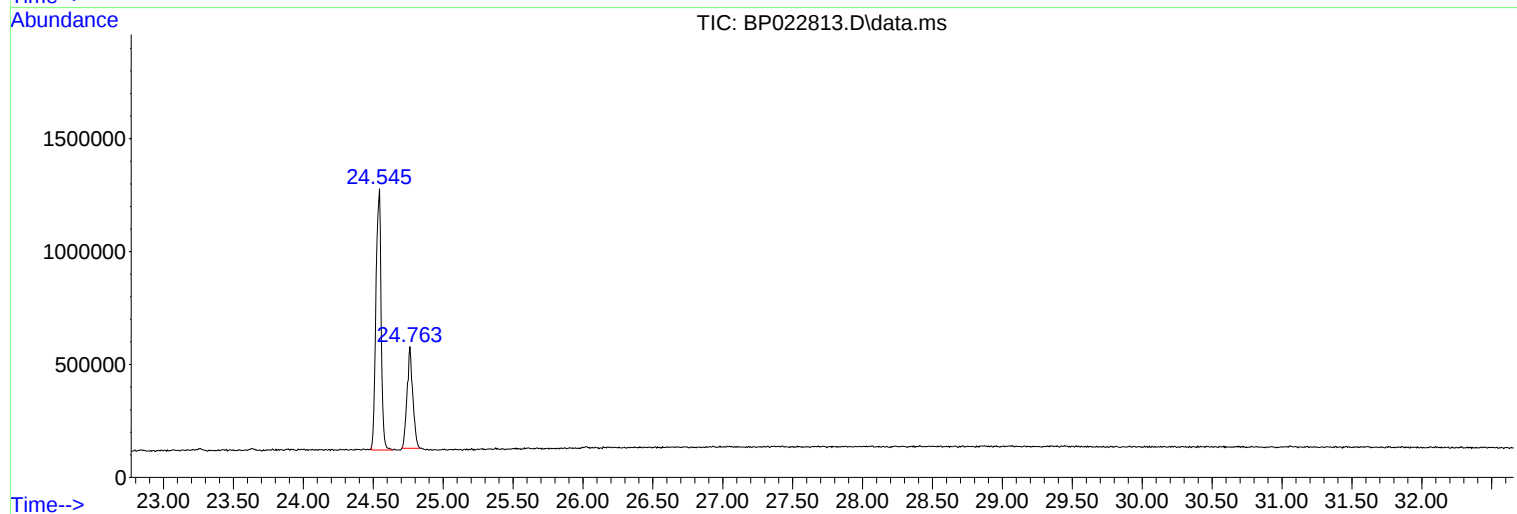
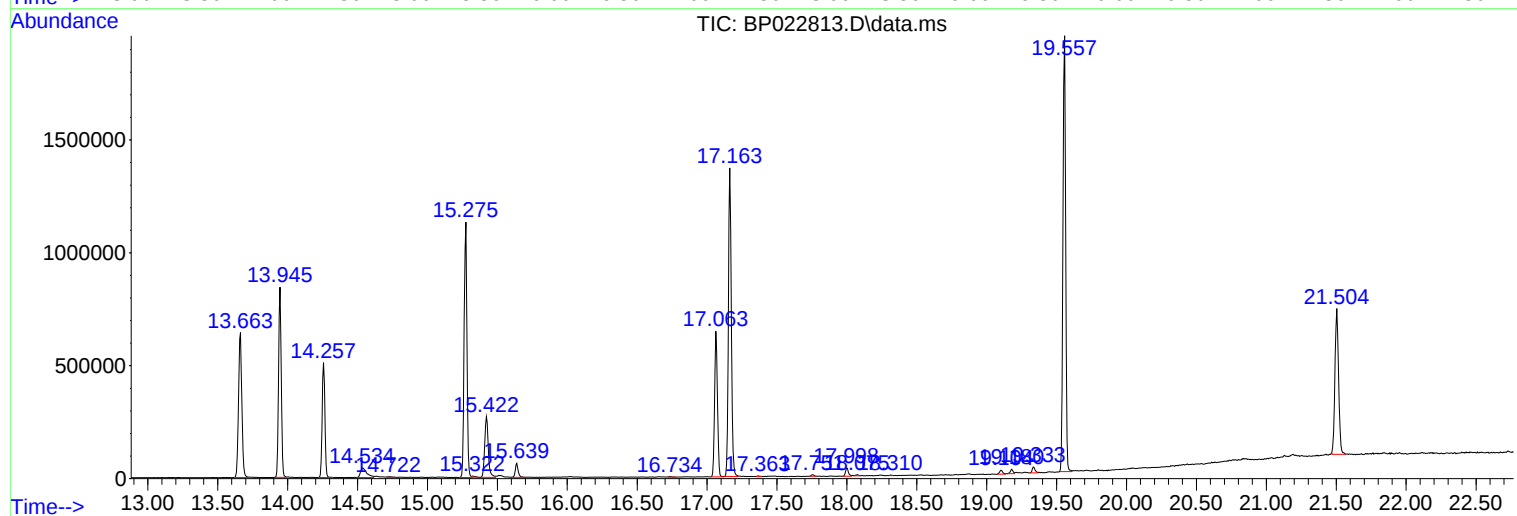
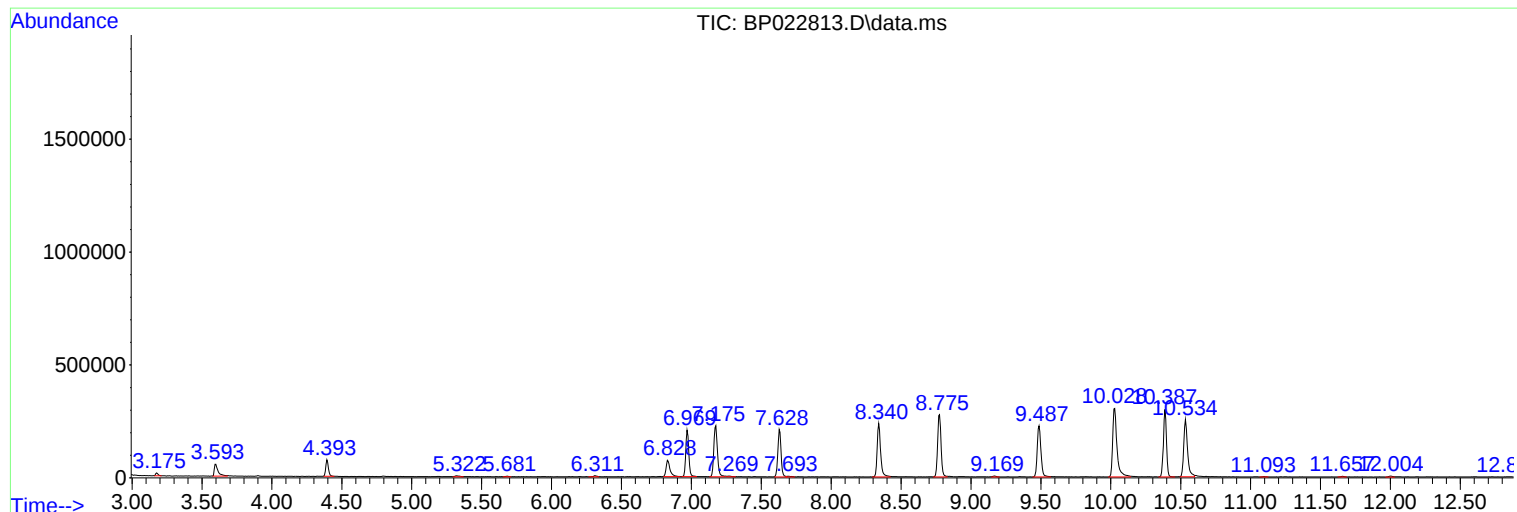
Sum of corrected areas: 20447935

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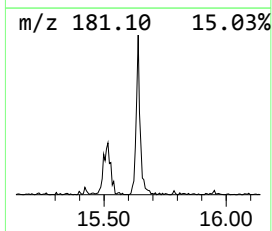
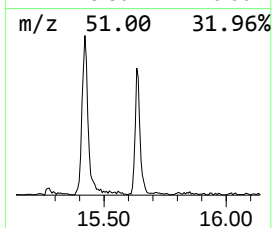
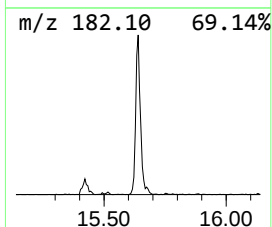
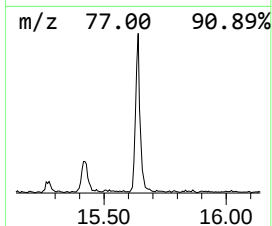
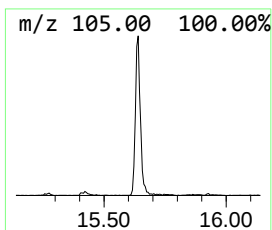
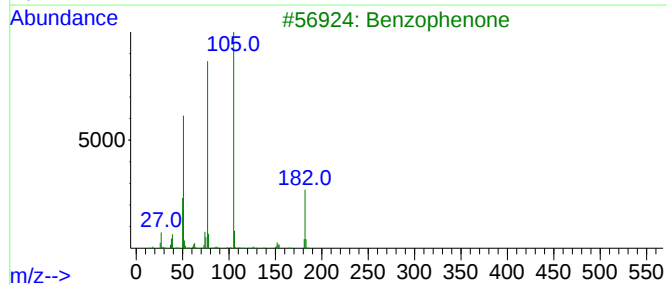
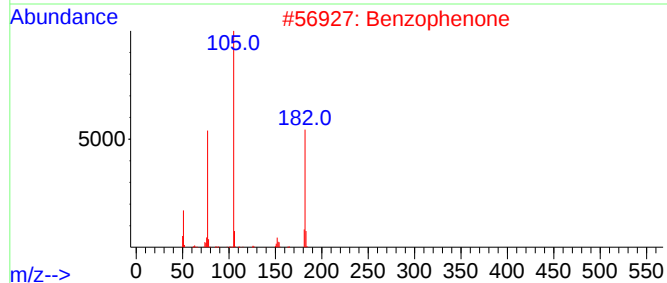
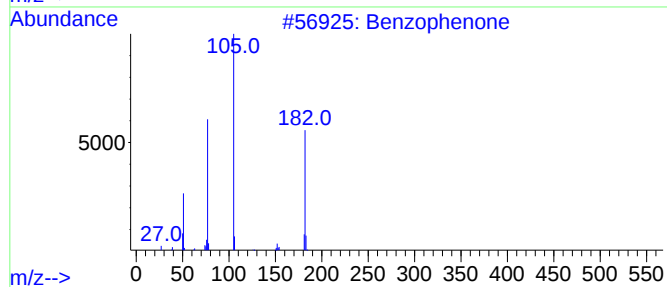
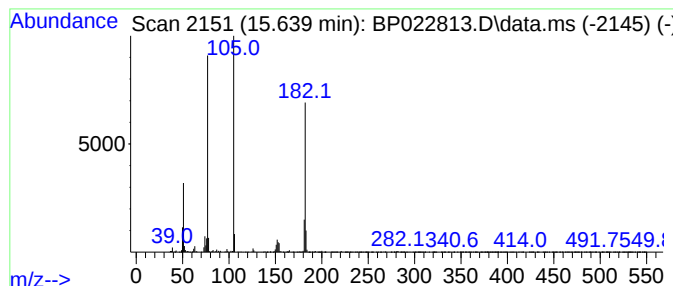
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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.
15.640	2.52 ng/ul	85810	Acenaphthene-d10	14.257

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	94
2			Benzophenone	182	C13H10O	000119-61-9	93
3			Benzophenone	182	C13H10O	000119-61-9	89
4			Benzophenone	182	C13H10O	000119-61-9	87
5			Benzophenone	182	C13H10O	000119-61-9	87



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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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.640	2.5	ng/u1	85810	3	14.257	682095	20.0