

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041223\
 Data File : BP014699.D
 Acq On : 13 Apr 2023 06:55
 Operator : CG/JU
 Sample : 02254-06
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AA8

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: BP014699.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.322	20	23	26	rBV4	13555	18441	0.48%	0.057%
2	3.358	26	29	32	rVV	29809	34978	0.91%	0.108%
3	3.393	32	35	43	rVB	38628	49132	1.27%	0.152%
4	3.811	105	106	116	rBV	51282	93584	2.43%	0.290%
5	4.105	153	156	162	rVB2	9448	13098	0.34%	0.041%
6	5.034	310	314	319	rBV8	3162	5744	0.15%	0.018%
7	6.052	484	487	495	rVB9	5955	11146	0.29%	0.035%
8	6.305	526	530	535	rBV3	9703	13581	0.35%	0.042%
9	7.146	666	673	691	rBV	86330	214293	5.56%	0.664%
10	7.299	693	699	715	rVV	441212	730898	18.95%	2.265%
11	7.499	726	733	751	rBV	429034	766518	19.87%	2.375%
12	7.969	806	813	823	rBV	404084	675608	17.52%	2.093%
13	8.081	828	832	842	rVB4	18244	33394	0.87%	0.103%
14	8.693	928	936	954	rBV	298293	586314	15.20%	1.817%
15	8.975	981	984	989	rBV7	3228	4724	0.12%	0.015%
16	9.146	1005	1013	1028	rBV	575100	1008286	26.14%	3.124%
17	9.505	1070	1074	1081	rVB9	4859	9043	0.23%	0.028%
18	9.869	1129	1136	1151	rBV	484905	864275	22.41%	2.678%
19	10.416	1221	1229	1251	rBV	498449	1096966	28.44%	3.399%
20	10.787	1284	1292	1300	rBV	582100	989814	25.66%	3.067%
21	10.940	1310	1318	1346	rBV	313622	784553	20.34%	2.431%
22	12.398	1557	1566	1574	rBV7	8942	21734	0.56%	0.067%
23	13.428	1730	1741	1744	rBV7	5945	13418	0.35%	0.042%
24	14.034	1837	1844	1856	rBV	1187323	1785440	46.29%	5.532%
25	14.222	1874	1876	1883	rVB7	2281	4276	0.11%	0.013%
26	14.316	1885	1892	1906	rVV	1370472	2029713	52.62%	6.289%
27	14.622	1936	1944	1956	rBV	856214	1241757	32.19%	3.847%
28	14.916	1987	1994	1997	rBV6	17525	39342	1.02%	0.122%
29	15.622	2106	2114	2125	rBV	1661615	2576116	66.79%	7.981%
30	15.751	2130	2136	2151	rVV	540211	914926	23.72%	2.835%
31	15.863	2152	2155	2162	rVB8	12772	23274	0.60%	0.072%
32	15.987	2171	2176	2190	rVB	159471	256748	6.66%	0.795%
33	16.198	2206	2212	2213	rBV6	3501	4691	0.12%	0.015%
34	16.557	2268	2273	2279	rBV2	21623	33449	0.87%	0.104%
35	17.057	2355	2358	2361	rBV5	2698	4029	0.10%	0.012%
36	17.175	2376	2378	2385	rVB8	3938	5905	0.15%	0.018%

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37	17.386	2407	2414	2424	rBV	1010300	1441788	37.38%	4.467%
38	17.486	2424	2431	2449	rVV	2019727	2944492	76.34%	9.123%
39	18.075	2528	2531	2535	rBV5	3749	6230	0.16%	0.019%
40	18.269	2560	2564	2571	rBV7	14907	31596	0.82%	0.098%
41	18.528	2605	2608	2610	rBV4	4342	5351	0.14%	0.017%
42	19.298	2735	2739	2744	rBV4	24020	33030	0.86%	0.102%
43	19.369	2747	2751	2757	rVV	21409	35935	0.93%	0.111%
44	19.498	2770	2773	2777	rBV6	12575	17924	0.46%	0.056%
45	19.716	2804	2810	2821	rBV	2522845	3489481	90.47%	10.811%
46	21.480	3104	3110	3120	rBV	1018483	1636212	42.42%	5.069%
47	23.810	3499	3506	3521	rBV	1896289	3857105	100.00%	11.950%
48	23.974	3528	3534	3551	rVB2	818565	1817994	47.13%	5.633%

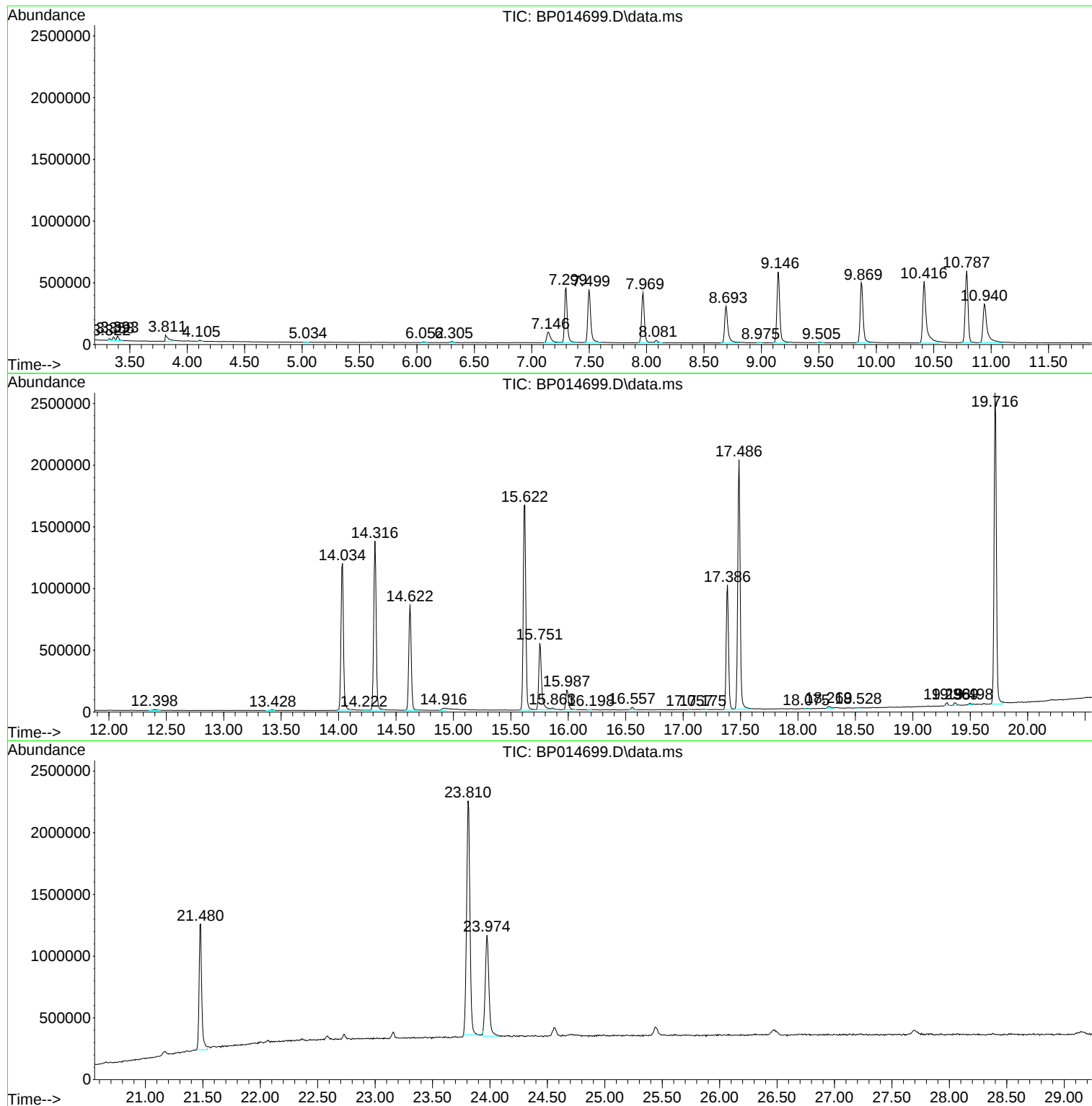
Sum of corrected areas: 32276346

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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



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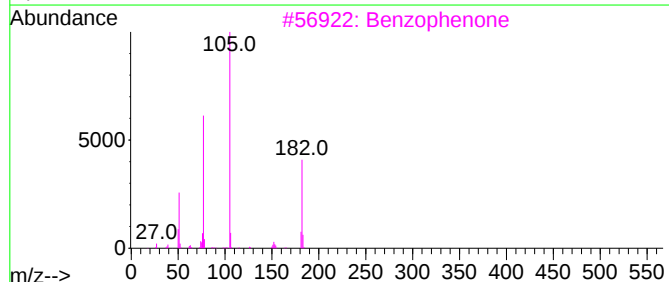
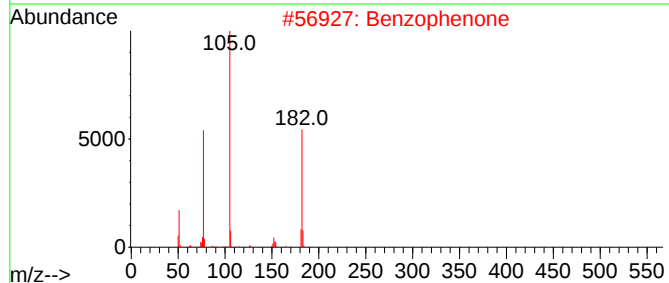
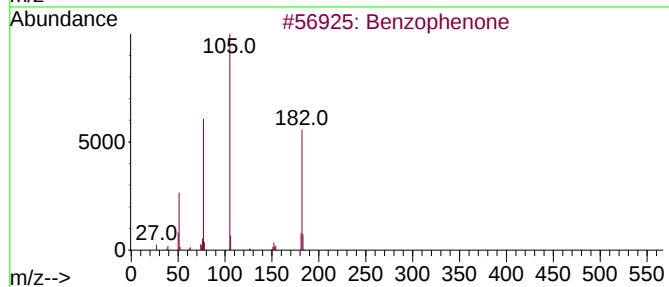
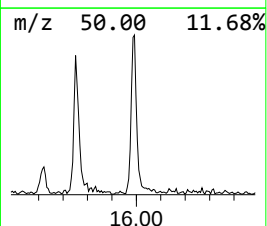
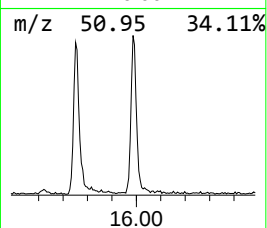
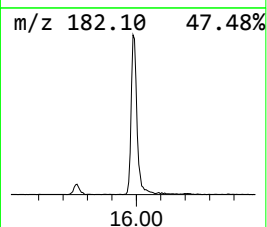
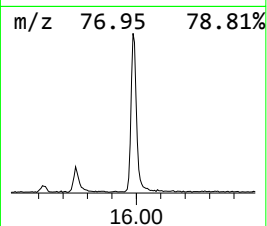
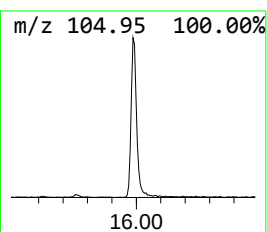
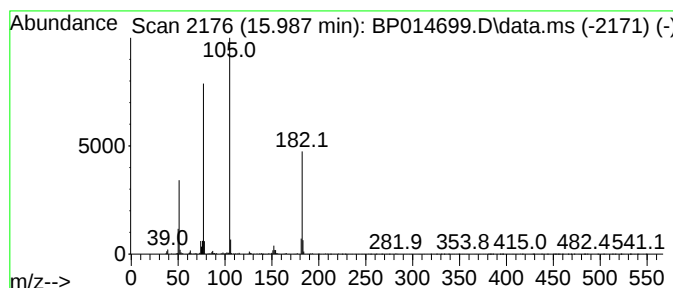
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
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TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.986	4.14 ng/ul	256748	Acenaphthene-d10	14.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	95
2			Benzophenone	182	C13H10O	000119-61-9	95
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 Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.986	4.1	ng/u1	256748	3	14.622	1241760	20.0