

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040323\
 Data File : BP014485.D
 Acq On : 03 Apr 2023 14:07
 Operator : CG/JU
 Sample : 02093-10DL 10X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 CB982DL

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: BP014485.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.417	36	39	53	rVB	93516	146255	5.68%	1.002%
2	3.911	122	123	127	rBV4	3290	5240	0.20%	0.036%
3	4.034	142	144	148	rVB5	3581	4555	0.18%	0.031%
4	4.140	157	162	165	rBV3	6990	10964	0.43%	0.075%
5	4.217	171	175	180	rBV	48489	62840	2.44%	0.431%
6	4.593	234	239	244	rBV7	8911	15493	0.60%	0.106%
7	4.881	286	288	293	rVB6	1842	2602	0.10%	0.018%
8	5.075	319	321	325	rVB5	2618	3064	0.12%	0.021%
9	6.958	637	641	643	rBV5	2517	3220	0.12%	0.022%
10	7.099	663	665	673	rVB9	3888	4642	0.18%	0.032%
11	7.205	679	683	692	rVB8	5580	13382	0.52%	0.092%
12	7.340	701	706	714	rBV	35317	59902	2.32%	0.411%
13	7.475	725	729	734	rBV8	2468	3967	0.15%	0.027%
14	7.546	734	741	755	rBV2	31515	71147	2.76%	0.488%
15	8.010	813	820	831	rBV	452063	777797	30.19%	5.331%
16	8.252	857	861	862	rBV3	3409	4115	0.16%	0.028%
17	8.499	899	903	906	rBV6	2617	4130	0.16%	0.028%
18	8.687	929	935	938	rVB8	2639	4160	0.16%	0.029%
19	8.752	938	946	960	rBV3	21380	68718	2.67%	0.471%
20	8.934	969	977	984	rBV3	5289	12282	0.48%	0.084%
21	9.022	984	992	1002	rBV	37929	103622	4.02%	0.710%
22	9.122	1008	1009	1016	rVB7	4844	6676	0.26%	0.046%
23	9.199	1016	1022	1029	rBV	37342	73516	2.85%	0.504%
24	9.381	1049	1053	1058	rVB7	7204	11449	0.44%	0.078%
25	9.499	1068	1073	1076	rBV7	1786	2774	0.11%	0.019%
26	9.528	1076	1078	1079	rBV2	3812	2939	0.11%	0.020%
27	9.640	1092	1097	1104	rBV4	26577	49815	1.93%	0.341%
28	9.734	1110	1113	1123	rBV4	4219	9426	0.37%	0.065%
29	9.922	1139	1145	1158	rBV3	30726	66089	2.56%	0.453%
30	10.034	1161	1164	1170	rVV8	4760	8873	0.34%	0.061%
31	10.099	1170	1175	1182	rVB8	4042	7473	0.29%	0.051%
32	10.369	1214	1221	1232	rBV8	8920	22047	0.86%	0.151%
33	10.487	1234	1241	1252	rBV3	29231	93723	3.64%	0.642%
34	10.757	1282	1287	1292	rVB9	3230	5898	0.23%	0.040%
35	10.834	1292	1300	1312	rBV	693658	1222211	47.43%	8.377%
36	10.916	1312	1314	1318	rVB5	3122	2970	0.12%	0.020%

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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
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37	10.957	1318	1321	1322	rBV3	4876	5363	0.21%	0.037%
38	11.004	1325	1329	1339	rVV4	17302	44386	1.72%	0.304%
39	11.093	1339	1344	1348	rVV7	5327	8665	0.34%	0.059%
40	11.551	1419	1422	1424	rBV4	2539	3102	0.12%	0.021%

41	11.681	1442	1444	1451	rVB7	5994	8183	0.32%	0.056%
42	11.746	1451	1455	1458	rBV6	2114	3416	0.13%	0.023%
43	11.804	1461	1465	1467	rBV5	4053	6178	0.24%	0.042%
44	11.863	1473	1475	1479	rBV5	2687	3674	0.14%	0.025%
45	12.046	1504	1506	1513	rVB8	4738	6194	0.24%	0.042%

46	12.175	1525	1528	1535	rBV9	3932	9136	0.35%	0.063%
47	12.281	1542	1546	1549	rVB6	2718	4335	0.17%	0.030%
48	12.369	1554	1561	1567	rBV6	13005	35977	1.40%	0.247%
49	12.463	1575	1577	1579	rBV3	4192	4551	0.18%	0.031%
50	12.557	1589	1593	1595	rVB5	2960	3504	0.14%	0.024%

51	12.840	1633	1641	1646	rBV9	8770	20386	0.79%	0.140%
52	13.045	1667	1676	1682	rBV9	6723	21637	0.84%	0.148%
53	13.145	1691	1693	1696	rBV4	2213	3081	0.12%	0.021%
54	13.216	1702	1705	1709	rBV6	2789	4329	0.17%	0.030%
55	13.469	1740	1748	1750	rBV9	4515	11240	0.44%	0.077%

56	13.545	1757	1761	1765	rVB5	10567	17183	0.67%	0.118%
57	13.681	1782	1784	1785	rBV2	4420	3530	0.14%	0.024%
58	13.710	1787	1789	1793	rVV5	3896	6432	0.25%	0.044%
59	14.075	1845	1851	1858	rVB	119303	187335	7.27%	1.284%
60	14.357	1894	1899	1907	rBV2	127201	202596	7.86%	1.389%

61	14.475	1914	1919	1929	rBV	30472	59901	2.32%	0.411%
62	14.587	1933	1938	1942	rBV4	20990	30753	1.19%	0.211%
63	14.663	1944	1951	1963	rBV	1177385	1737027	67.41%	11.906%
64	14.987	2004	2006	2009	rBV4	4360	5100	0.20%	0.035%
65	15.175	2036	2038	2041	rBV4	3721	3822	0.15%	0.026%

66	15.663	2115	2121	2129	rBV	181957	283493	11.00%	1.943%
67	15.798	2140	2144	2155	rVB4	35911	74358	2.89%	0.510%
68	16.186	2205	2210	2220	rVB	37712	63500	2.46%	0.435%
69	16.704	2294	2298	2302	rBV4	18206	32321	1.25%	0.222%
70	17.281	2394	2396	2401	rBV6	8468	13135	0.51%	0.090%

71	17.428	2415	2421	2431	rBV	1541614	2267716	88.01%	15.543%
72	17.528	2433	2438	2452	rVB	220640	340353	13.21%	2.333%
73	18.292	2565	2568	2577	rBV8	29974	47713	1.85%	0.327%
74	19.757	2812	2817	2822	rBV	310599	402105	15.61%	2.756%
75	21.516	3111	3116	3124	rVB	1964797	2576747	100.00%	17.661%

76	23.868	3512	3516	3523	rVB	328480	585841	22.74%	4.015%
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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

77 24.033 3538 3544 3554 rVB2 1181814 2473738 96.00% 16.955%

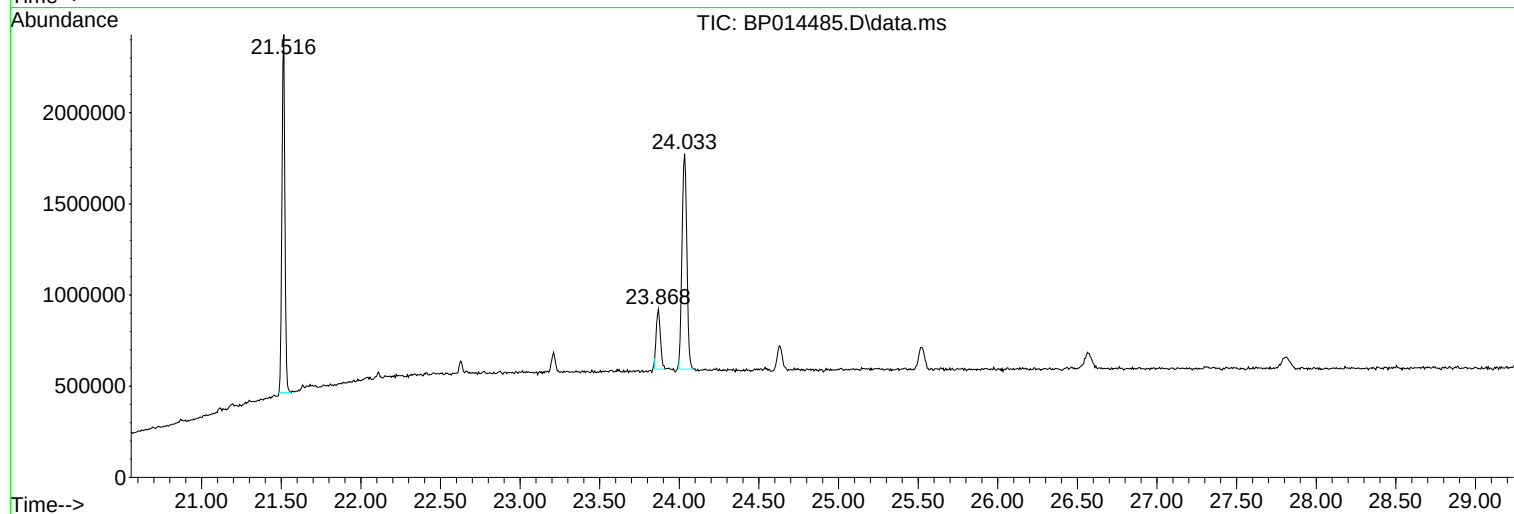
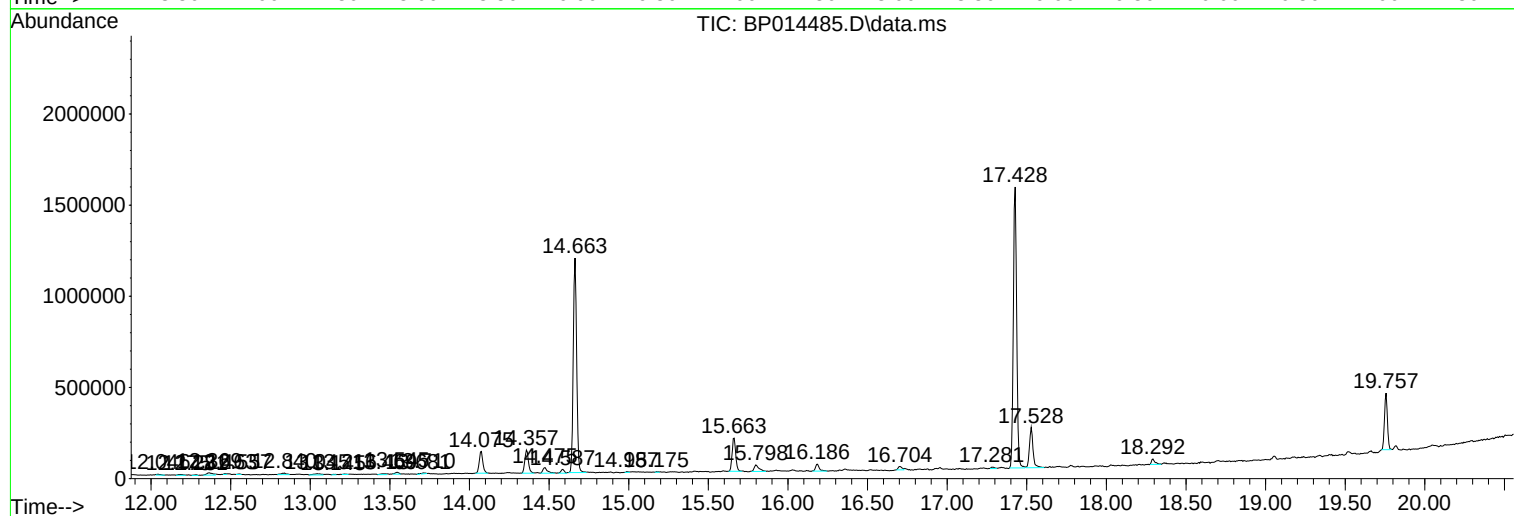
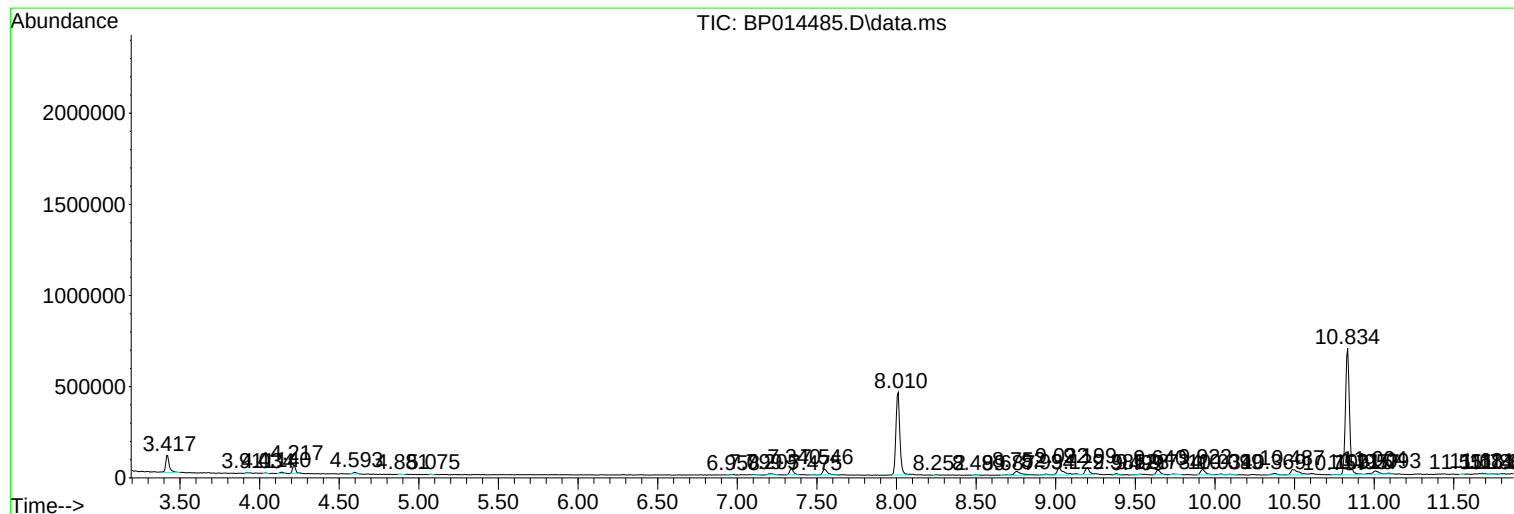
Sum of corrected areas: 14590012

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Instrument :
BNA_P
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CB982DL

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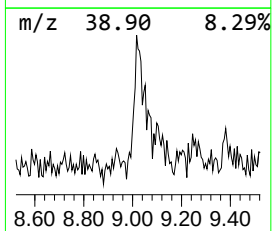
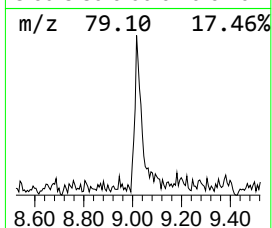
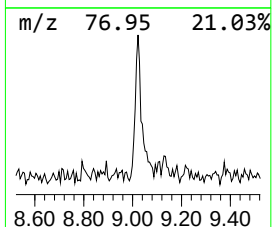
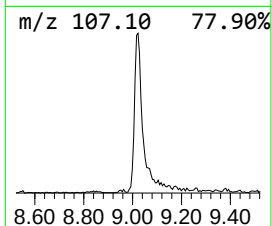
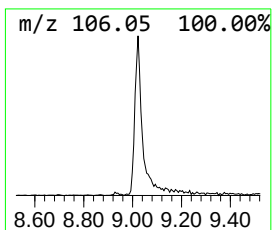
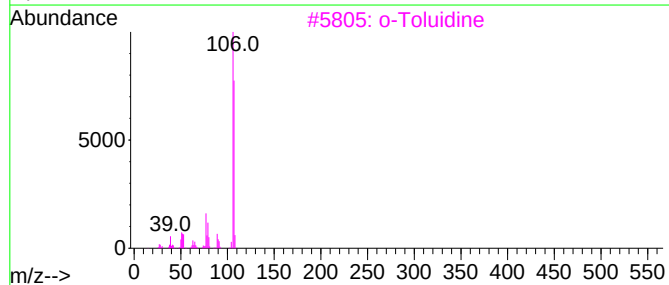
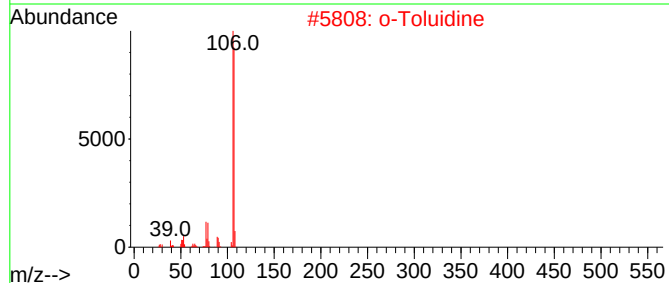
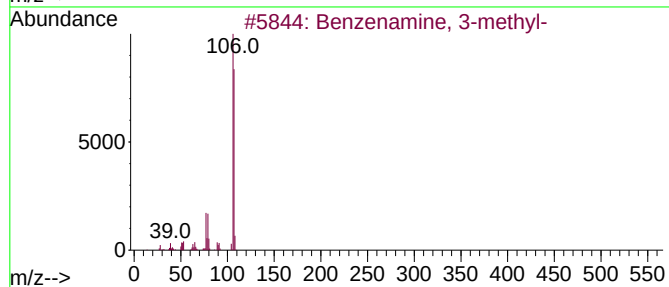
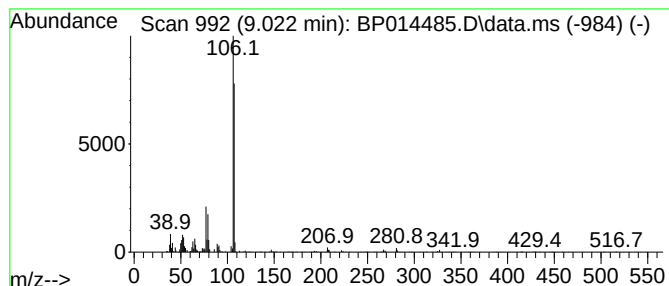
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzenamine, 3-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.022	2.66 ng/ul	103622	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	93
2			o-Toluidine	107	C7H9N	000095-53-4	90
3			o-Toluidine	107	C7H9N	000095-53-4	90
4			Benzenamine, 3-methyl-	107	C7H9N	000108-44-1	90
5			p-Aminotoluene	107	C7H9N	000106-49-0	90



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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
Benzenamine, 3-...	9.022	2.7	ng/u1	103622	1	8.010	777797	20.0