

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110923\
 Data File : BP018010.D
 Acq On : 10 Nov 2023 03:28
 Operator : MA/JU
 Sample : 05091-03
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCK12

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-BP110923.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018010.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.858	128	131	137	rVB2	10630	15932	2.20%	0.295%
2	4.910	306	310	315	rVB2	12123	17380	2.40%	0.321%
3	5.905	477	479	483	rBV4	1748	2367	0.33%	0.044%
4	5.981	490	492	495	rBV3	1405	1195	0.16%	0.022%
5	6.340	551	553	556	rBV4	1594	1358	0.19%	0.025%
6	7.046	667	673	690	rBV2	34057	107059	14.76%	1.979%
7	7.216	697	702	710	rBV	33298	59717	8.23%	1.104%
8	7.393	726	732	741	rBV2	47974	90360	12.46%	1.671%
9	7.610	767	769	773	rVB6	2055	1802	0.25%	0.033%
10	7.652	773	776	777	rBV3	1697	1691	0.23%	0.031%
11	7.816	802	804	806	rVB3	1466	1278	0.18%	0.024%
12	7.863	806	812	826	rBV	150257	259449	35.77%	4.797%
13	8.134	854	858	859	rBV4	1368	1866	0.26%	0.035%
14	8.263	877	880	881	rBV3	1202	1382	0.19%	0.026%
15	8.510	920	922	926	rVB5	1529	1783	0.25%	0.033%
16	8.557	926	930	932	rBV5	1278	1261	0.17%	0.023%
17	8.604	932	938	946	rBV4	21751	52104	7.18%	0.963%
18	8.734	953	960	971	rBV	60891	121690	16.78%	2.250%
19	9.081	1013	1019	1031	rBV	41582	88116	12.15%	1.629%
20	9.687	1121	1122	1125	rBV3	1117	1249	0.17%	0.023%
21	9.793	1132	1140	1150	rBV2	32879	76895	10.60%	1.422%
22	10.175	1200	1205	1214	rBV8	5357	10998	1.52%	0.203%
23	10.334	1227	1232	1243	rBV2	31341	79809	11.00%	1.476%
24	10.587	1272	1275	1280	rBV6	1401	2096	0.29%	0.039%
25	10.687	1285	1292	1304	rBV	191296	338072	46.61%	6.251%
26	10.887	1320	1326	1338	rBV3	20145	58172	8.02%	1.076%
27	11.204	1377	1380	1384	rBV5	1183	1770	0.24%	0.033%
28	11.257	1384	1389	1391	rBV5	1382	2436	0.34%	0.045%
29	11.304	1395	1397	1399	rVB3	1796	1373	0.19%	0.025%
30	11.469	1423	1425	1430	rVB6	1518	1558	0.21%	0.029%
31	11.610	1447	1449	1452	rBV4	1394	1667	0.23%	0.031%
32	11.669	1458	1459	1462	rBV3	1471	1442	0.20%	0.027%
33	11.740	1466	1471	1472	rVV5	1301	1910	0.26%	0.035%
34	11.822	1483	1485	1487	rVB3	1392	1195	0.16%	0.022%
35	11.845	1487	1489	1494	rBV6	1074	1778	0.25%	0.033%
36	12.028	1516	1520	1522	rVB4	1129	1236	0.17%	0.023%

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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-BP110923.MA.M
 Title : SVOA CALIBRATION

37	12.187	1543	1547	1550	rVB5	1870	1581	0.22%	0.029%
38	12.292	1561	1565	1569	rBV7	1202	2005	0.28%	0.037%
39	12.487	1594	1598	1601	rVB6	1267	1455	0.20%	0.027%
40	12.528	1601	1605	1607	rBV5	1437	1622	0.22%	0.030%
41	12.557	1607	1610	1612	rBV4	962	1190	0.16%	0.022%
42	12.669	1628	1629	1631	rBV2	1526	1375	0.19%	0.025%
43	12.734	1638	1640	1645	rVB6	1352	1983	0.27%	0.037%
44	12.957	1675	1678	1681	rBV5	1440	1765	0.24%	0.033%
45	13.098	1701	1702	1706	rVB4	1394	1248	0.17%	0.023%
46	13.322	1737	1740	1743	rBV4	813	1222	0.17%	0.023%
47	13.381	1743	1750	1756	rVV4	8365	16899	2.33%	0.312%
48	13.757	1810	1814	1817	rBV5	1089	1428	0.20%	0.026%
49	13.845	1826	1829	1831	rBV4	1087	1333	0.18%	0.025%
50	13.963	1843	1849	1856	rBV	108315	177347	24.45%	3.279%
51	14.128	1872	1877	1878	rVB5	1259	1664	0.23%	0.031%
52	14.239	1887	1896	1907	rBV2	119987	194426	26.80%	3.595%
53	14.416	1922	1926	1927	rVB4	1176	1240	0.17%	0.023%
54	14.539	1941	1947	1955	rBV	298721	438292	60.43%	8.104%
55	14.628	1960	1962	1968	rVB7	2817	2704	0.37%	0.050%
56	14.728	1977	1979	1982	rBV4	1385	1808	0.25%	0.033%
57	14.857	1995	2001	2005	rBV9	7413	16855	2.32%	0.312%
58	15.128	2045	2047	2051	rBV5	1974	1997	0.28%	0.037%
59	15.281	2069	2073	2075	rBV4	2178	2472	0.34%	0.046%
60	15.369	2083	2088	2095	rBV	7296	11720	1.62%	0.217%
61	15.539	2112	2117	2125	rBV	161809	257377	35.48%	4.759%
62	15.704	2140	2145	2153	rBV	27980	52126	7.19%	0.964%
63	15.986	2191	2193	2196	rBV3	1139	1273	0.18%	0.024%
64	16.192	2226	2228	2232	rBV5	1188	1408	0.19%	0.026%
65	16.316	2245	2249	2253	rBV7	1813	2625	0.36%	0.049%
66	16.416	2264	2266	2271	rVB4	2883	3567	0.49%	0.066%
67	16.833	2336	2337	2338	rBV	2876	1926	0.27%	0.036%
68	16.975	2359	2361	2363	rBV3	1431	1549	0.21%	0.029%
69	17.316	2413	2419	2425	rBV	398617	558531	77.00%	10.327%
70	17.422	2431	2437	2452	rVB	184138	287701	39.66%	5.319%
71	18.163	2559	2563	2572	rBV5	27252	41621	5.74%	0.770%
72	18.992	2701	2704	2709	rBV2	11640	14604	2.01%	0.270%
73	19.339	2760	2763	2771	rVB	42600	61938	8.54%	1.145%
74	19.669	2814	2819	2823	rBV	311538	433066	59.71%	8.007%
75	21.439	3116	3120	3125	rBV	487866	666876	91.94%	12.330%
76	23.951	3542	3547	3563	rVB	345778	725341	100.00%	13.411%

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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-BP110923.MA.M
Title : SVOA CALIBRATION

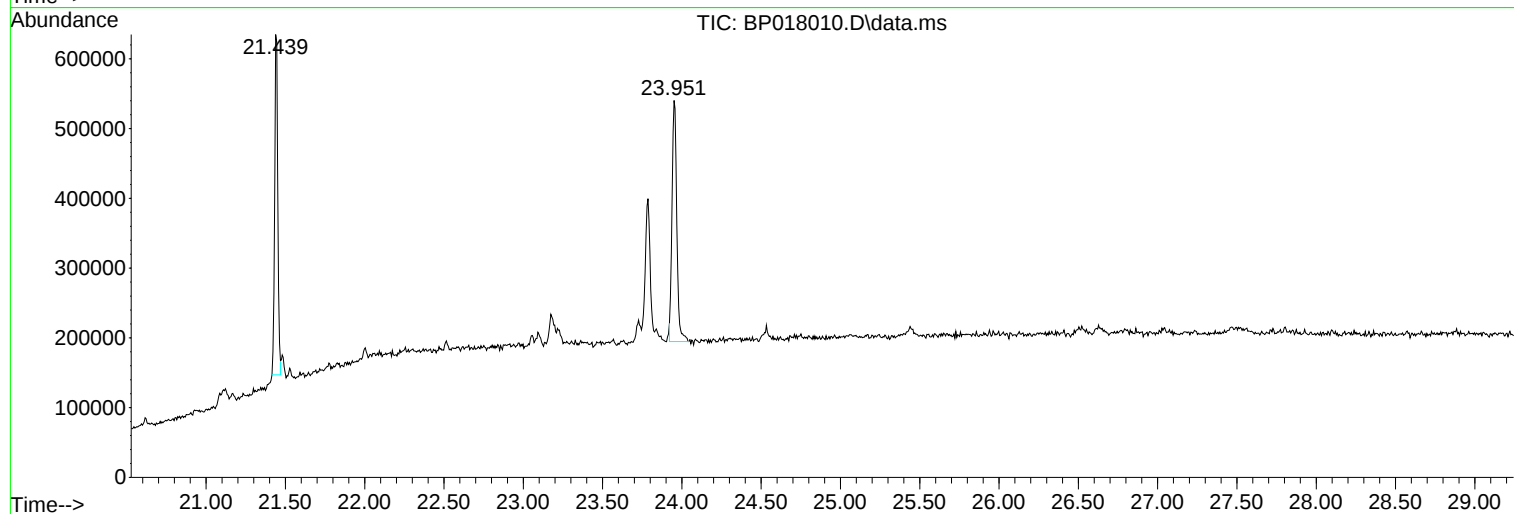
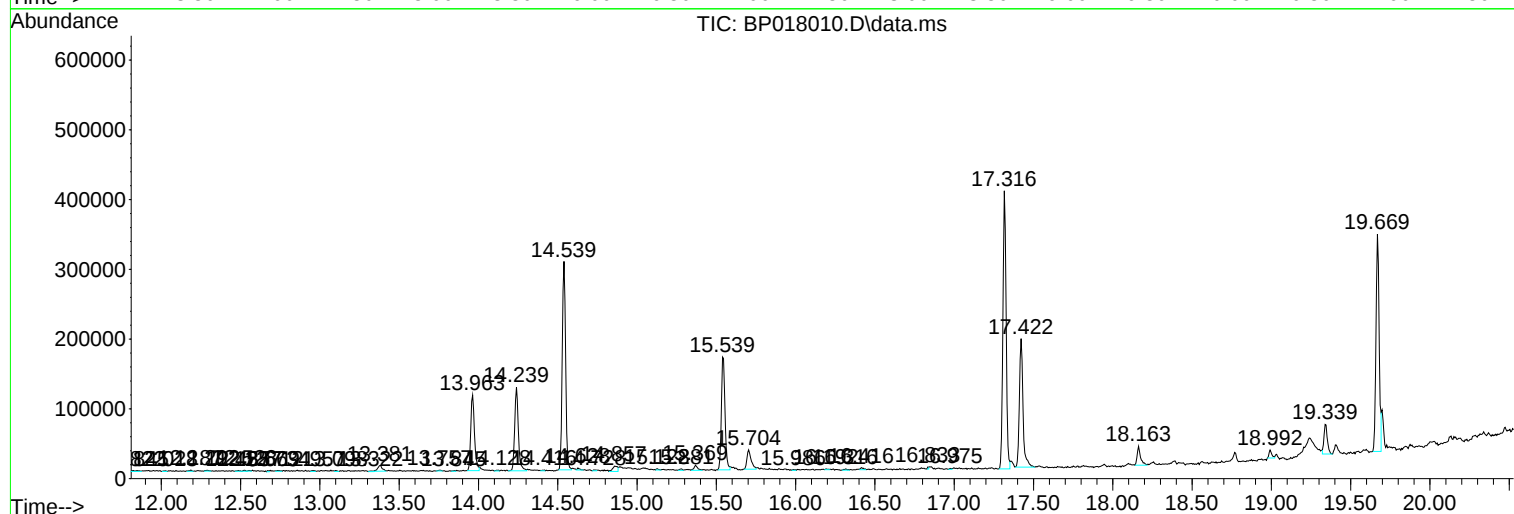
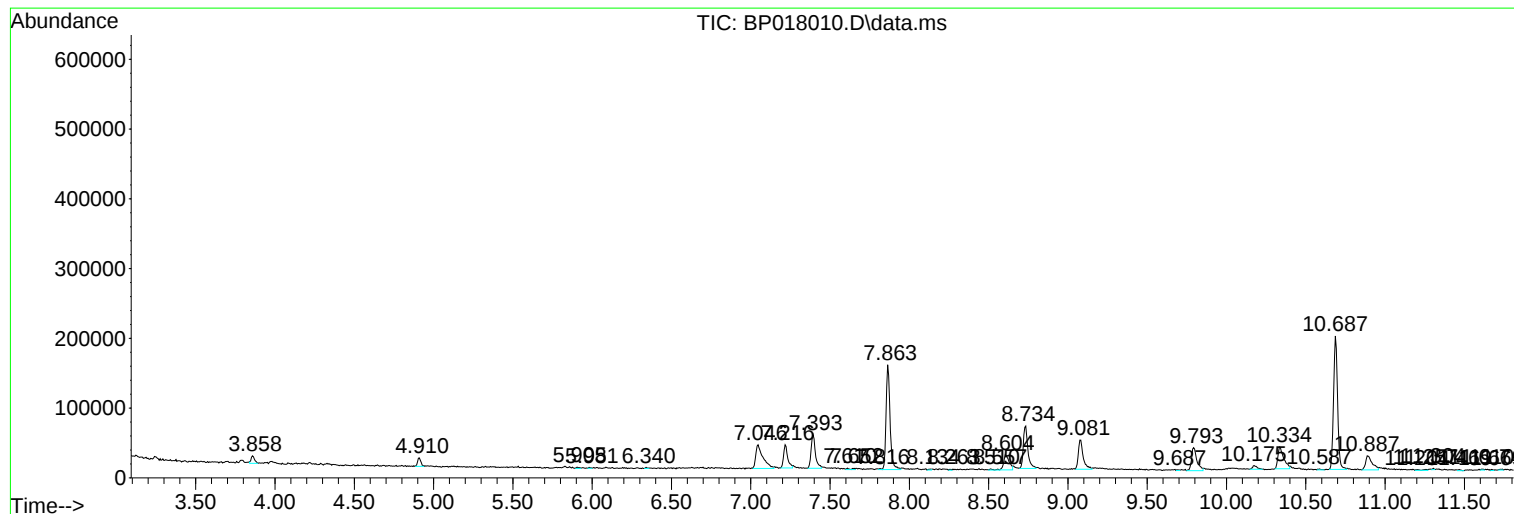
Sum of corrected areas: 5408606

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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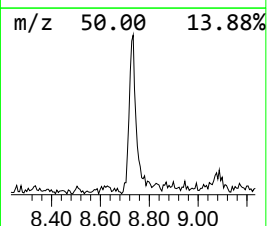
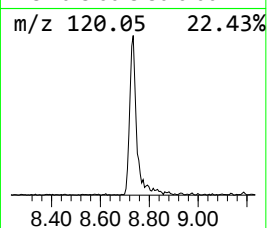
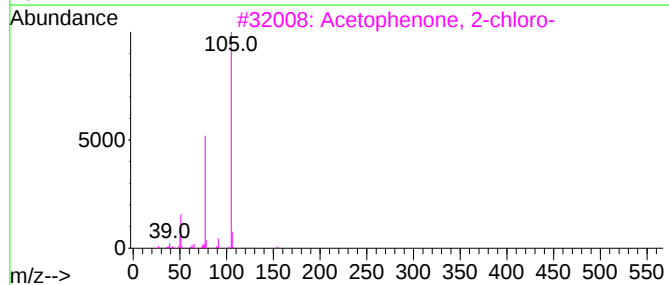
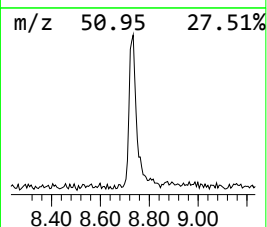
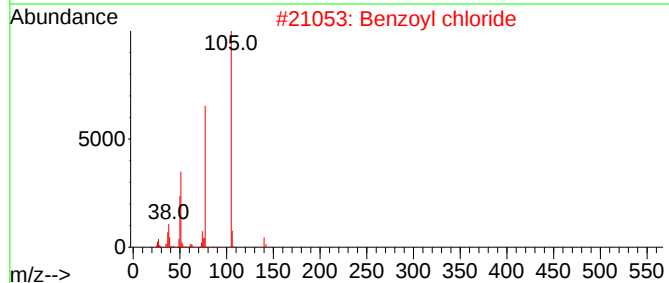
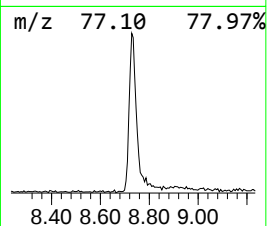
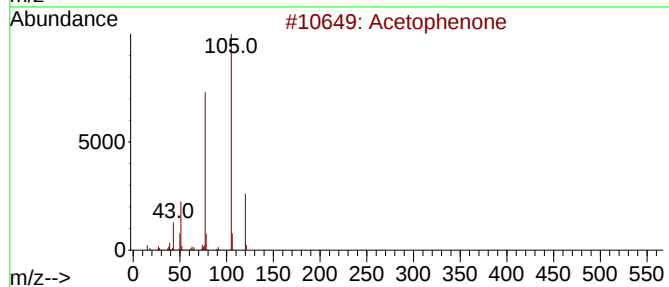
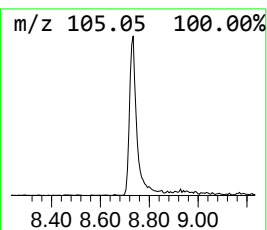
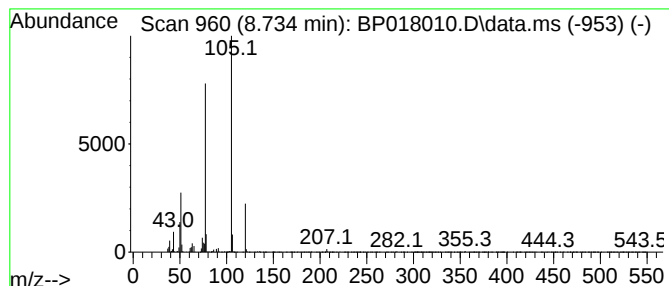
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110923.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Aceto phenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.734	9.38 ng/ul	121690	1,4-Dichlorobenzene-d4	7.863

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetophenone	120	C8H8O	000098-86-2	90
2			Benzoyl chloride	140	C7H5ClO	000098-88-4	87
3			Acetophenone, 2-chloro-	154	C8H7ClO	000532-27-4	80
4			Benzoyl bromide	184	C7H5BrO	000618-32-6	80
5			Ethanone, 2,2-dihydroxy-1-phenyl-	152	C8H8O3	001075-06-5	80



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Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Aceto phenone	8.734	9.4	ng/u1	121690	1	7.863	259449	20.0