

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM101119\
 Data File : BM023145.D
 Acq On : 12 Oct 2019 00:19
 Operator : JU
 Sample : PB123768BL
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK68

Integration Parameters: LSCINT.P

Integrator: RTE
 Smoothing : OFF
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 0.1 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
 Title : SVOA CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.252	42	45	52	rBV	48740	64110	1.60%	0.156%
2	3.975	165	168	175	rVB	10041	13799	0.34%	0.034%
3	6.939	665	672	688	rBV	809056	1266801	31.58%	3.078%
4	7.098	693	699	719	rBV	590594	986888	24.61%	2.397%
5	7.304	727	734	744	rVB	840953	1369275	34.14%	3.326%
6	7.769	806	813	821	rBV	709311	1123362	28.01%	2.729%
7	8.475	926	933	950	rBV	968587	1545979	38.55%	3.756%
8	8.922	1002	1009	1026	rBV	726038	1198873	29.89%	2.912%
9	9.639	1124	1131	1144	rBV	620845	1005619	25.07%	2.443%
10	10.180	1215	1223	1237	rBV	922046	1585242	39.52%	3.851%
11	10.557	1279	1287	1296	rVB	909721	1549843	38.64%	3.765%
12	10.692	1303	1310	1325	rBV	847989	1530360	38.16%	3.718%
13	12.145	1553	1557	1564	rVB	12412	19089	0.48%	0.046%
14	13.816	1834	1841	1854	rBV	1404037	1939794	48.36%	4.712%
15	14.098	1882	1889	1902	rBV	1729110	2485306	61.96%	6.038%
16	14.404	1935	1941	1952	rVB	1228006	1805882	45.03%	4.387%
17	14.610	1968	1976	1994	rBV	471139	812879	20.27%	1.975%
18	15.398	2103	2110	2122	rBV	2098332	2839631	70.80%	6.898%
19	15.515	2125	2130	2143	rBV	462978	650037	16.21%	1.579%
20	15.639	2146	2151	2157	rVB	19072	25862	0.64%	0.063%
21	15.957	2194	2205	2212	rBV3	17827	32667	0.81%	0.079%
22	16.180	2240	2243	2246	rVB3	7226	8060	0.20%	0.020%
23	16.633	2315	2320	2325	rBV4	9445	15679	0.39%	0.038%
24	16.939	2368	2372	2375	rVB	3198	4978	0.12%	0.012%
25	17.145	2401	2407	2414	rBV2	1308772	1863411	46.46%	4.527%
26	17.245	2417	2424	2437	rVV	2111061	2918250	72.76%	7.089%
27	17.398	2445	2450	2456	rBV6	4571	8473	0.21%	0.021%
28	19.174	2748	2752	2756	rBV	21764	28418	0.71%	0.069%
29	19.427	2792	2795	2798	rBV	19130	22773	0.57%	0.055%
30	19.539	2808	2814	2821	rBV	2601633	3434651	85.63%	8.344%
31	21.033	3064	3068	3074	rBV	230256	301838	7.53%	0.733%
32	21.327	3113	3118	3125	rBV	1754499	2266671	56.51%	5.507%
33	23.438	3470	3477	3491	rVB	2112153	4010824	100.00%	9.744%
34	23.580	3495	3501	3511	rVB	1237399	2427962	60.54%	5.898%

LSC Area Percent Report

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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

Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
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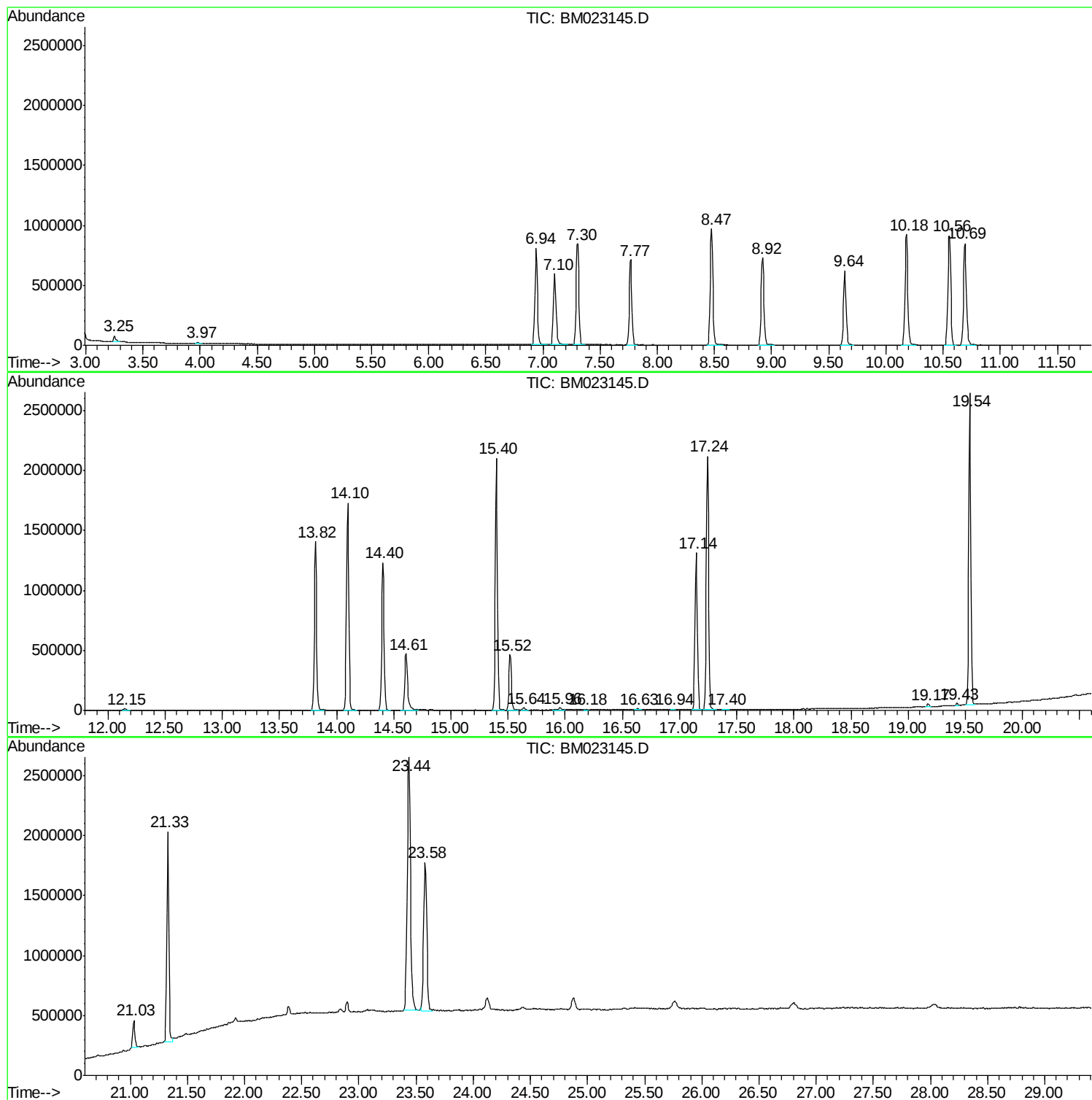
Sum of corrected areas: 41163286

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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM092719MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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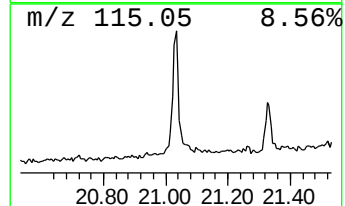
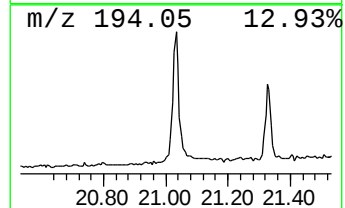
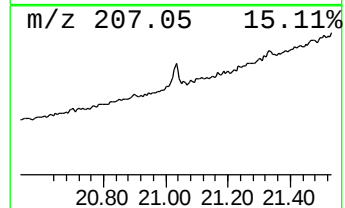
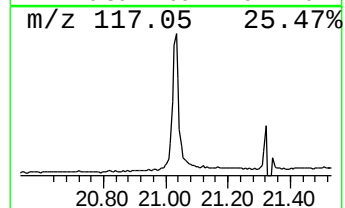
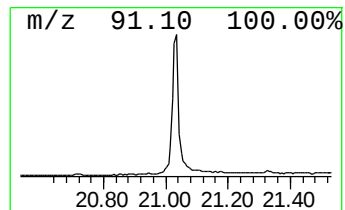
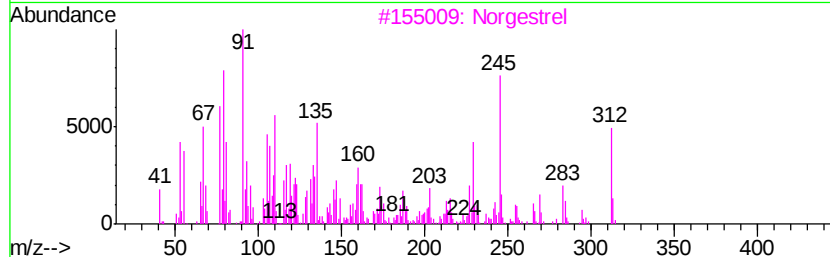
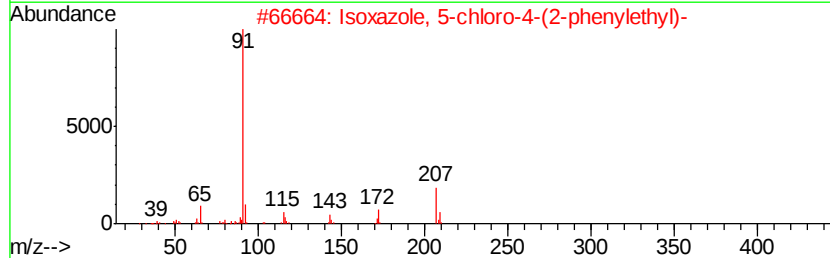
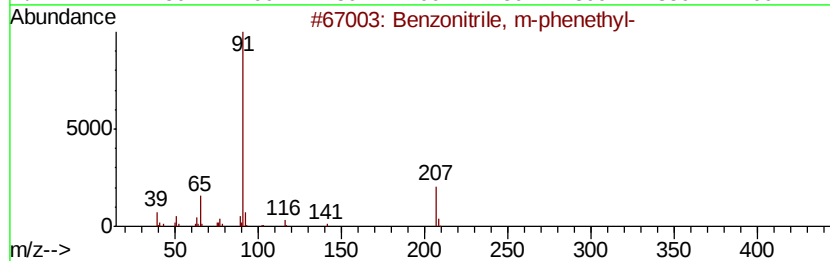
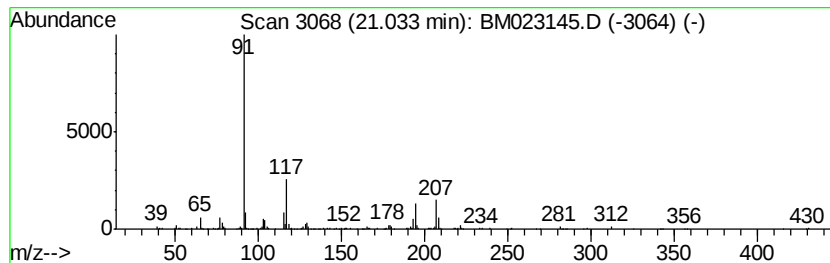
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Peak Number 1 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.03	2.66 ng/ul	301838	Chrysene-d12	21.33

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzonitrile, m-phenethyl-	207	C15H13N	034176-91-5	16
2		Isoxazole, 5-chloro-4-(2-phenylethyl)-	207	C11H10ClNO	1000362-09-0	16
3		Norgestrel	312	C21H28O2	006533-00-2	11
4		4-Benzylloxy-3-methoxy-2-nitrobenzyl	303	C15H13NO6	003584-32-5	10
5		Benzyl phenyl sulfone	232	C13H12O2S	003112-88-7	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	21.03	2.7	ng/ul	301838	5	21.33	2266670	20.0