

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122415\
 Data File : BG020475.D
 Acq On : 24 Dec 2015 16:29
 Operator : UM/SJ
 Sample : PB87437BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK37

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:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG122415.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	2.949	18	19	24	rVB	1341	1067	0.14%	0.017%
2	3.096	39	44	51	rBV	11913	21720	2.76%	0.337%
3	3.172	54	57	64	rBV2	3758	5382	0.68%	0.084%
4	3.449	100	104	109	rBV	4112	5558	0.71%	0.086%
5	4.342	253	256	259	rBV3	548	827	0.10%	0.013%
6	5.740	491	494	499	rVB2	592	951	0.12%	0.015%
7	7.232	741	748	759	rBV	73264	129798	16.47%	2.015%
8	7.397	770	776	785	rBV	61466	102704	13.03%	1.594%
9	7.608	806	812	819	rBV	76641	133465	16.94%	2.072%
10	8.078	886	892	901	rVB	52616	86257	10.95%	1.339%
11	8.784	1006	1012	1023	rBV	99589	172966	21.95%	2.685%
12	9.248	1084	1091	1100	rBV	79017	141158	17.91%	2.191%
13	9.970	1207	1214	1223	rBV	71822	124362	15.78%	1.930%
14	10.517	1300	1307	1319	rBV	100480	180971	22.97%	2.809%
15	10.893	1364	1371	1380	rBV	78023	133359	16.92%	2.070%
16	11.028	1387	1394	1403	rBV2	97825	182700	23.19%	2.836%
17	12.485	1638	1642	1648	rBB	778	1597	0.20%	0.025%
18	14.113	1912	1919	1929	rBV	235529	333534	42.33%	5.177%
19	14.406	1963	1969	1985	rVB	239428	374912	47.58%	5.820%
20	14.712	2015	2021	2030	rBB2	135872	216995	27.54%	3.368%
21	14.912	2049	2055	2065	rBV	86008	136098	17.27%	2.113%
22	15.705	2183	2190	2198	rBV	333647	502334	63.75%	7.798%
23	15.816	2204	2209	2217	rBV	102743	152667	19.37%	2.370%
24	15.940	2227	2230	2235	rVB	2885	3335	0.42%	0.052%
25	17.456	2482	2488	2495	rBV2	201505	300448	38.13%	4.664%
26	17.556	2498	2505	2516	rVV2	387804	589986	74.87%	9.158%
27	17.826	2549	2551	2555	rVB2	551	813	0.10%	0.013%
28	18.549	2671	2674	2679	rVB3	538	824	0.10%	0.013%
29	19.477	2827	2832	2837	rBV2	4510	6191	0.79%	0.096%
30	19.577	2847	2849	2854	rBV2	567	835	0.11%	0.013%
31	19.730	2871	2875	2879	rVB	3597	4491	0.57%	0.070%
32	19.841	2887	2894	2902	rBV	541781	758630	96.28%	11.776%
33	19.994	2918	2920	2924	rBV3	819	795	0.10%	0.012%
34	20.070	2931	2933	2935	rBV2	1186	916	0.12%	0.014%

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35	20.106	2935	2939	2940	rBV2	990	1377	0.17%	0.021%
36	20.393	2985	2988	2989	rBV2	1486	1059	0.13%	0.016%
37	20.534	3010	3012	3014	rVB2	1275	1140	0.14%	0.018%
38	20.564	3014	3017	3019	rBV2	1600	1743	0.22%	0.027%
39	20.599	3019	3023	3025	rVV4	1427	1613	0.20%	0.025%
40	20.910	3074	3076	3080	rBV4	1978	1791	0.23%	0.028%
41	21.386	3154	3157	3159	rBV4	1772	1884	0.24%	0.029%
42	21.486	3172	3174	3177	rVB4	1311	1052	0.13%	0.016%
43	21.733	3209	3216	3224	rBV	246096	413949	52.53%	6.426%
44	21.909	3244	3246	3250	rBV4	1530	2133	0.27%	0.033%
45	24.007	3601	3603	3604	rBV2	2024	1678	0.21%	0.026%
46	24.771	3721	3733	3746	rBV	281107	787962	100.00%	12.231%
47	24.994	3761	3771	3787	rVB	143549	404847	51.38%	6.284%
48	25.200	3804	3806	3811	rBV5	1766	2553	0.32%	0.040%
49	26.099	3956	3959	3960	rBV2	2203	2641	0.34%	0.041%
50	29.260	4495	4497	4499	rVB2	1522	1005	0.13%	0.016%
51	29.312	4504	4506	4507	rBV2	1490	988	0.13%	0.015%
52	29.418	4522	4524	4526	rBV3	1536	1457	0.18%	0.023%
53	30.963	4785	4787	4788	rBV	1573	981	0.12%	0.015%
54	31.727	4915	4917	4919	rVB2	1910	1599	0.20%	0.025%

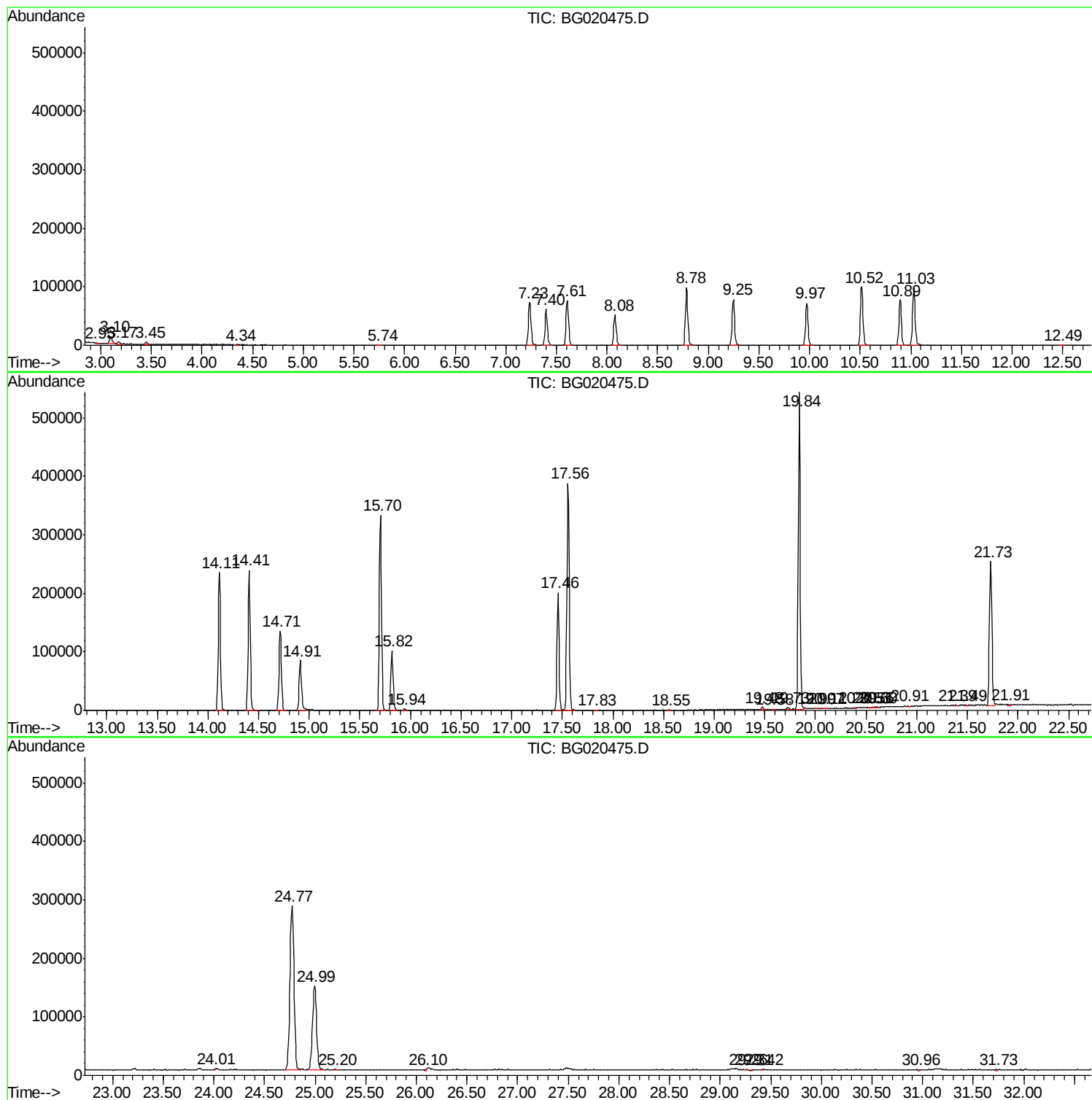
Sum of corrected areas: 6442098

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG122415.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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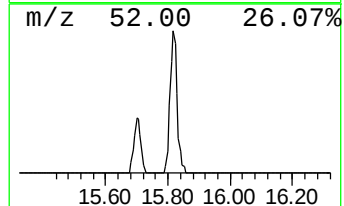
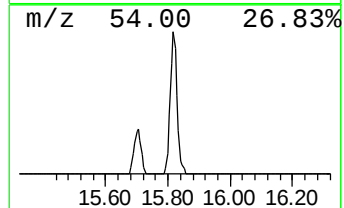
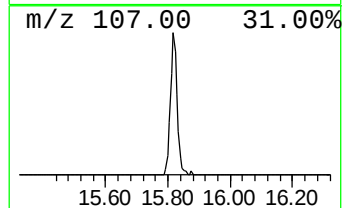
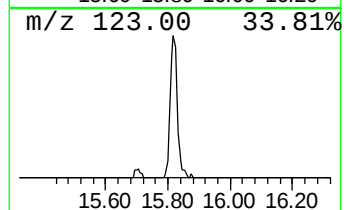
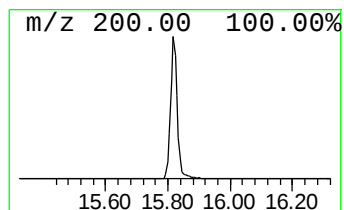
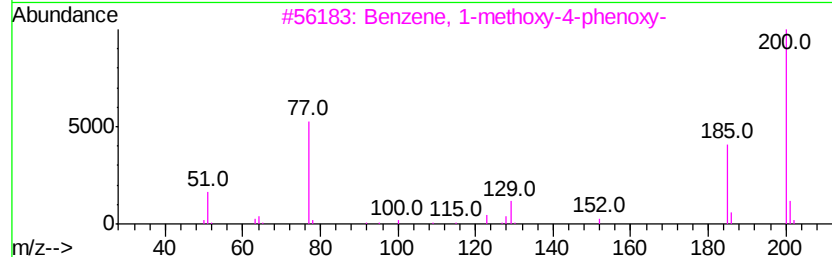
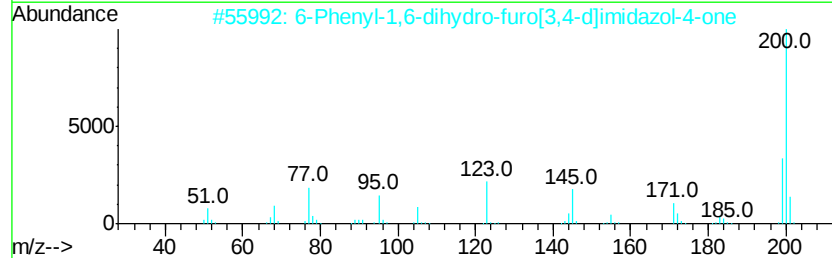
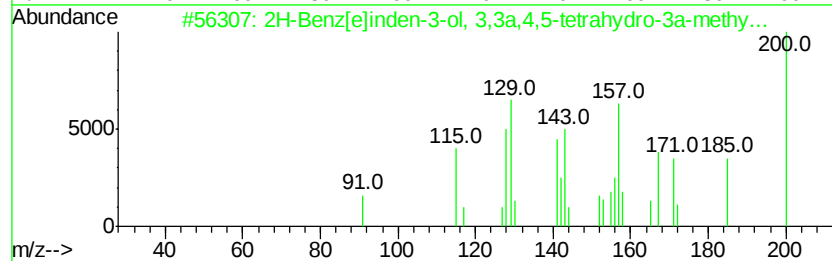
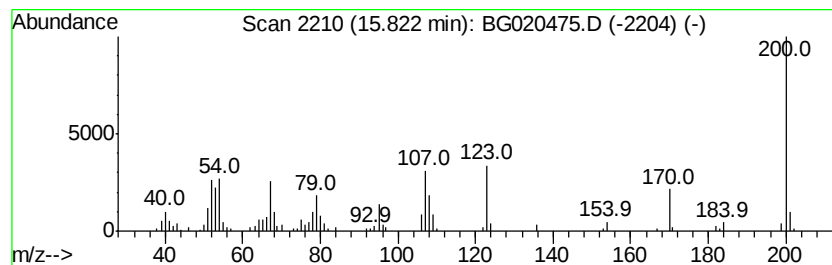
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Peak Number 2 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.82	14.07 ng/ul	152667	Acenaphthene-d10	14.71

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2H-Benz[e]inden-3-ol, 3,3a,4,5-t...	200	C14H16O	071805-91-9	35
2		6-Phenyl-1,6-dihydro-furo[3,4-d]...	200	C11H8N2O2	313263-12-6	25
3		Benzene, 1-methoxy-4-phenoxy-	200	C13H12O2	001655-69-2	22
4		Benzene, 1-methoxy-3-phenoxy-	200	C13H12O2	001655-68-1	22
5		1,6-Dioxaspiro[4.4]non-3-ene, 2-...	200	C13H12O2	016863-61-9	16



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	15.82	14.1	ng/ul	152667	3	14.71	216995	20.0