

Data Path : Z:\HPCHEM1\BNA G\DATA\BG112217\
 Data File : BG031199.D
 Acq On : 22 Nov 2017 21:11
 Operator : SJ/JU
 Sample : I6525-11
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 T71-MW5

Integration Parameters: rteint.p
 Integrator: RTE
 Smoothing : ON Filtering: 5
 Sampling : 1 Min Area: 3 % 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:\HPCHEM1\BNA_G\METHODS\8270-BG111017.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.694	402	409	416	rBV	485166	820144	12.71%	3.393%
2	5.753	416	419	429	rVB	31027	69032	1.07%	0.286%
3	7.298	675	682	697	rBV	300721	589145	9.13%	2.437%
4	7.668	738	745	759	rBV	889285	1594558	24.71%	6.597%
5	8.132	818	824	833	rBV	192048	343750	5.33%	1.422%
6	8.461	872	880	890	rBV	842317	1496029	23.19%	6.189%
7	9.301	1015	1023	1041	rBV	612125	1189900	18.44%	4.923%
8	10.952	1295	1304	1324	rBV	248282	536756	8.32%	2.221%
9	13.379	1710	1717	1729	rBV	2106701	3471113	53.80%	14.360%
10	14.760	1941	1952	1966	rBV2	453798	829957	12.86%	3.434%
11	16.240	2196	2204	2226	rBV	1678919	2729710	42.31%	11.293%
12	17.498	2410	2418	2432	rBV2	716927	1146817	17.77%	4.745%
13	20.089	2852	2859	2869	rBV	4733919	6451862	100.00%	26.692%
14	21.763	3138	3144	3161	rBV	853746	1498846	23.23%	6.201%
15	25.065	3694	3706	3721	rBV2	468592	1403697	21.76%	5.807%

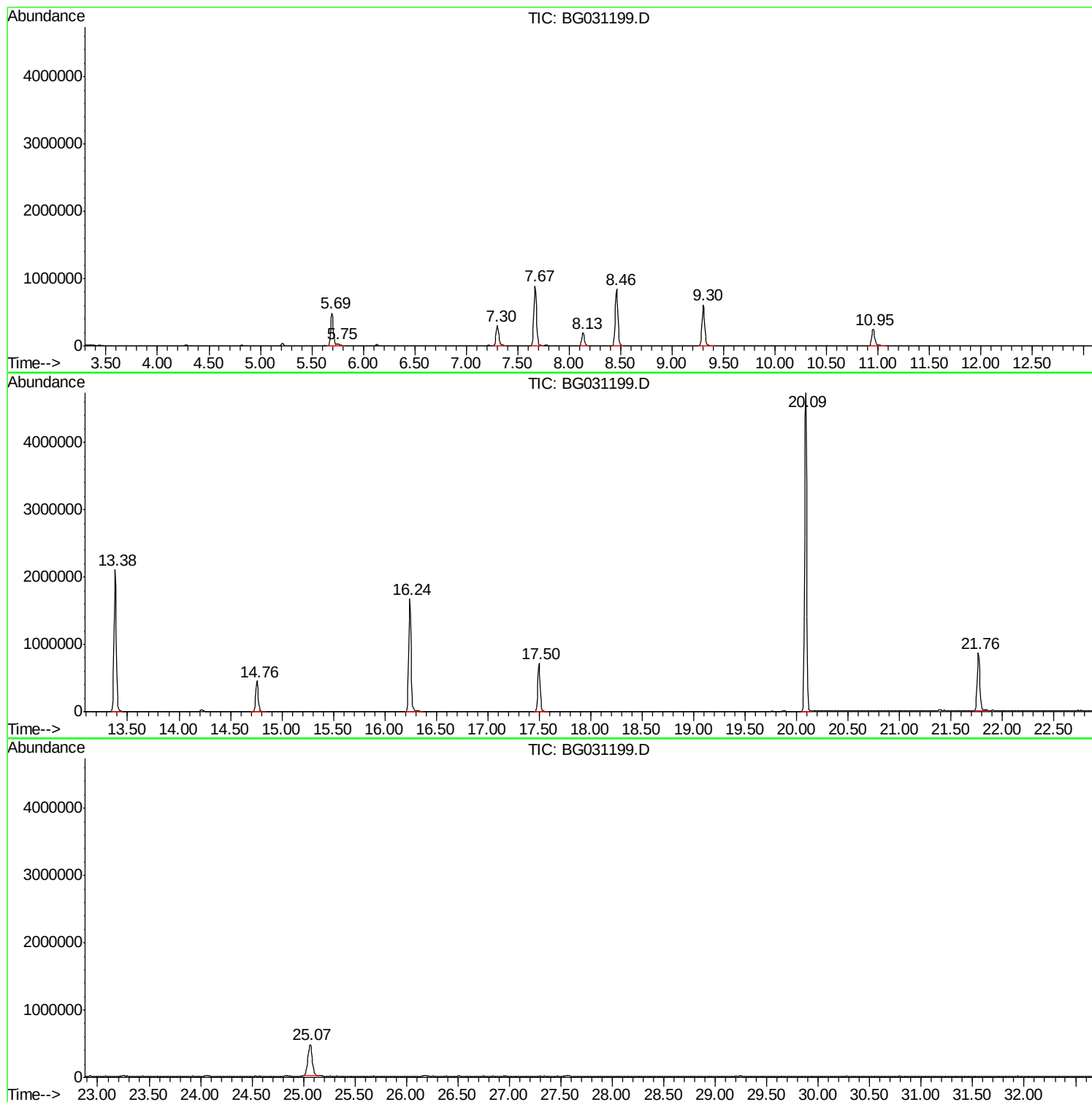
Sum of corrected areas: 24171316

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG111017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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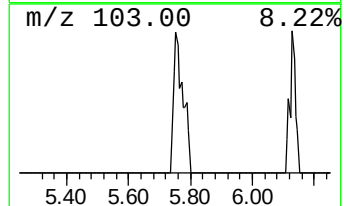
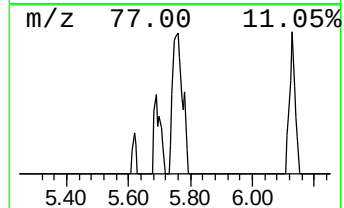
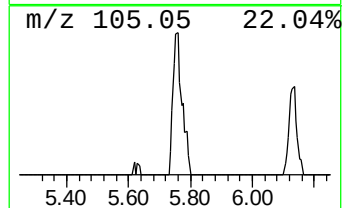
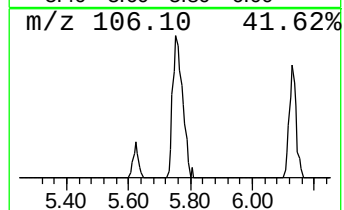
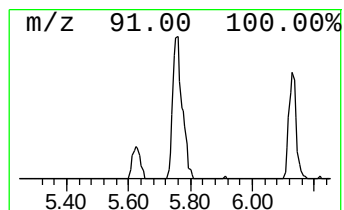
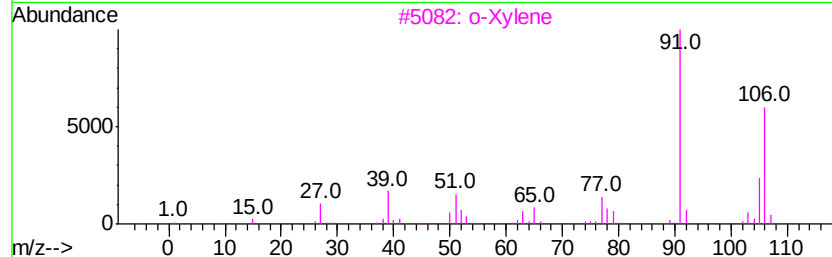
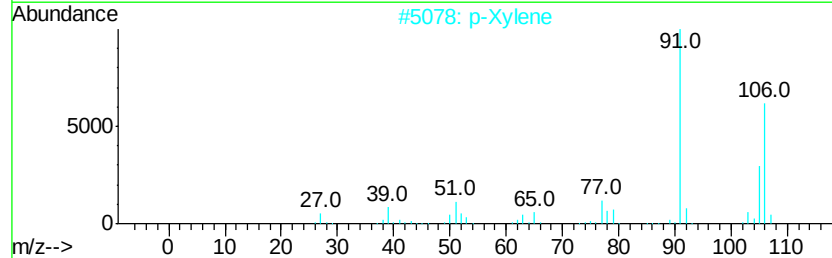
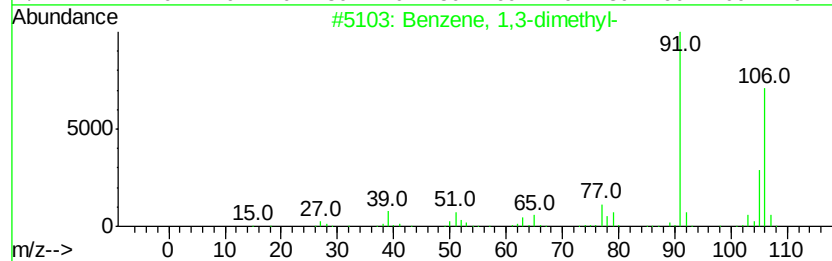
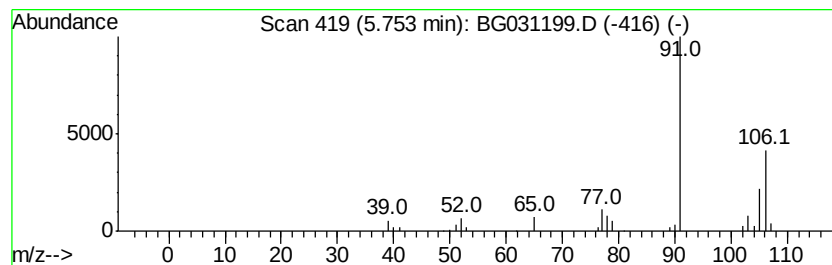
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Peak Number 1 Benzene, 1,3-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.75	4.02 ng	69032	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	91
2		p-Xylene	106	C8H10	000106-42-3	91
3		o-Xylene	106	C8H10	000095-47-6	91
4		Cyclopentene, 1-ethenyl-3-methyl...	106	C8H10	061142-07-2	72
5		Ethylbenzene	106	C8H10	000100-41-4	64



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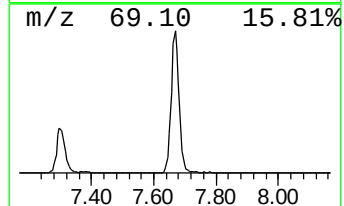
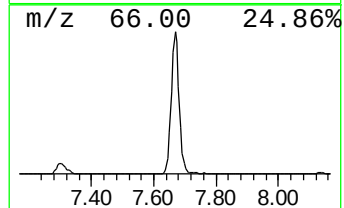
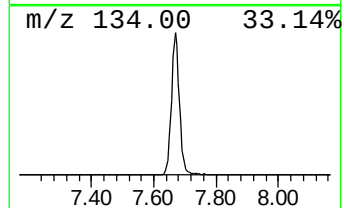
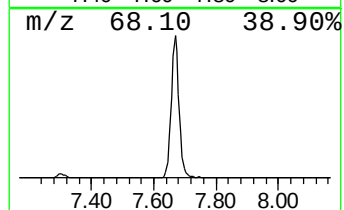
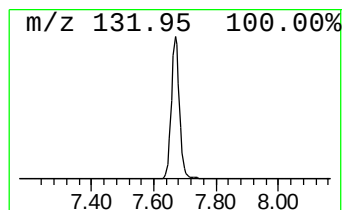
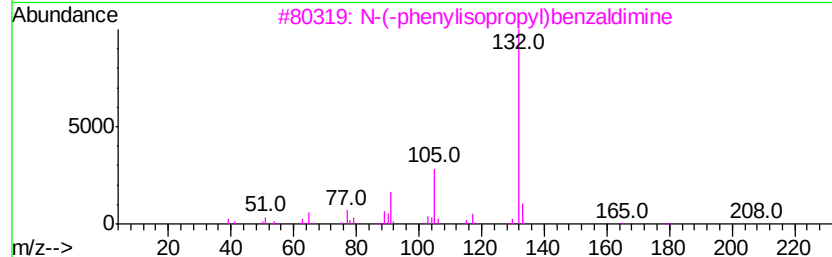
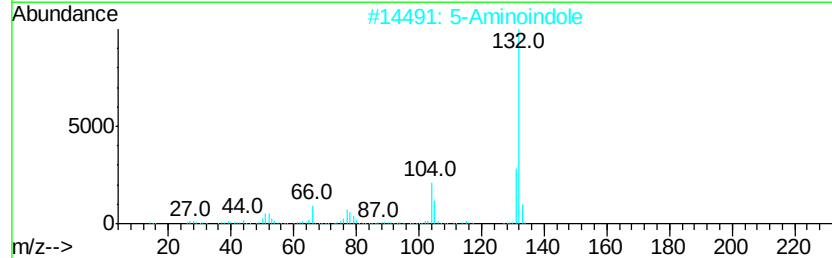
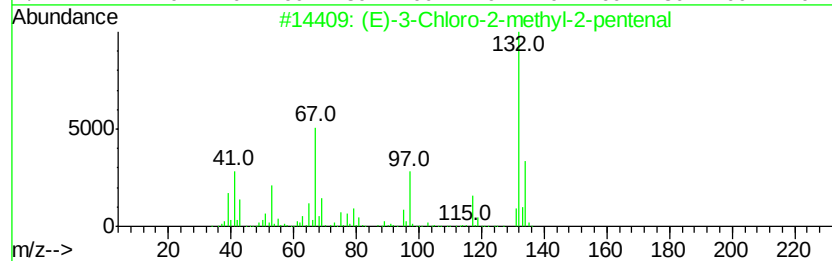
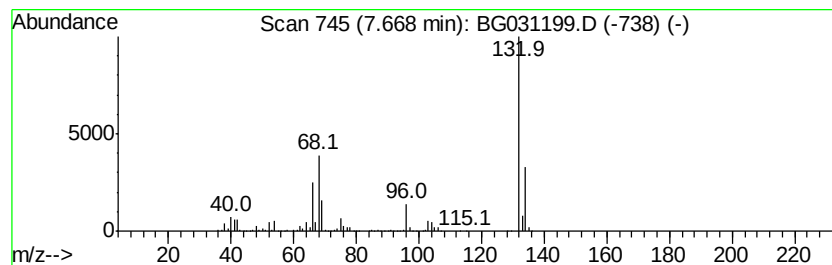
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Peak Number 2 unknown7.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	92.77 ng	1594560	1,4-Dichlorobenzene-d4	8.13

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		5-Aminoindole	132	C8H8N2	005192-03-0	12
3		N-(-phenylisopropyl)benzalimine	223	C16H17N	1000379-01-3	10
4		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-dime...	5.75	4.0	ng	69032	1	8.13	343750	20.0
unknown7.67	7.67	92.8	ng	1594560	1	8.13	343750	20.0