

Data Path : Z:\HPCHEM1\BNA G\Data\BG032517\
 Data File : BG026416.D
 Acq On : 24 Mar 2017 15:54
 Operator : SJ/MA
 Sample : PB97805BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB97805BL

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-BG032017.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	5.144	344	350	360	rBV	193024	296678	3.17%	0.647%
2	5.620	424	431	446	rBV	2203826	3505291	37.41%	7.650%
3	7.212	693	702	720	rBV	1980698	3387160	36.15%	7.392%
4	7.576	757	764	778	rBV	2122808	3753551	40.06%	8.192%
5	8.046	834	844	853	rBV	440252	760579	8.12%	1.660%
6	8.369	891	899	909	rBV	1774998	3075273	32.82%	6.712%
7	9.198	1033	1040	1055	rBV	1153884	2104609	22.46%	4.593%
8	10.843	1311	1320	1333	rBV	590953	1085277	11.58%	2.369%
9	13.275	1727	1734	1753	rBV	3623320	5789480	61.79%	12.635%
10	14.650	1961	1968	1977	rBV2	936106	1455866	15.54%	3.177%
11	16.131	2213	2220	2236	rBV	2717041	4215061	44.99%	9.199%
12	17.382	2426	2433	2447	rBV2	1148632	1833673	19.57%	4.002%
13	19.979	2869	2875	2891	rBV	7097046	9369522	100.00%	20.448%
14	21.624	3148	3155	3169	rBV	1565976	2599173	27.74%	5.673%
15	24.803	3686	3696	3710	rVB2	900730	2589104	27.63%	5.651%

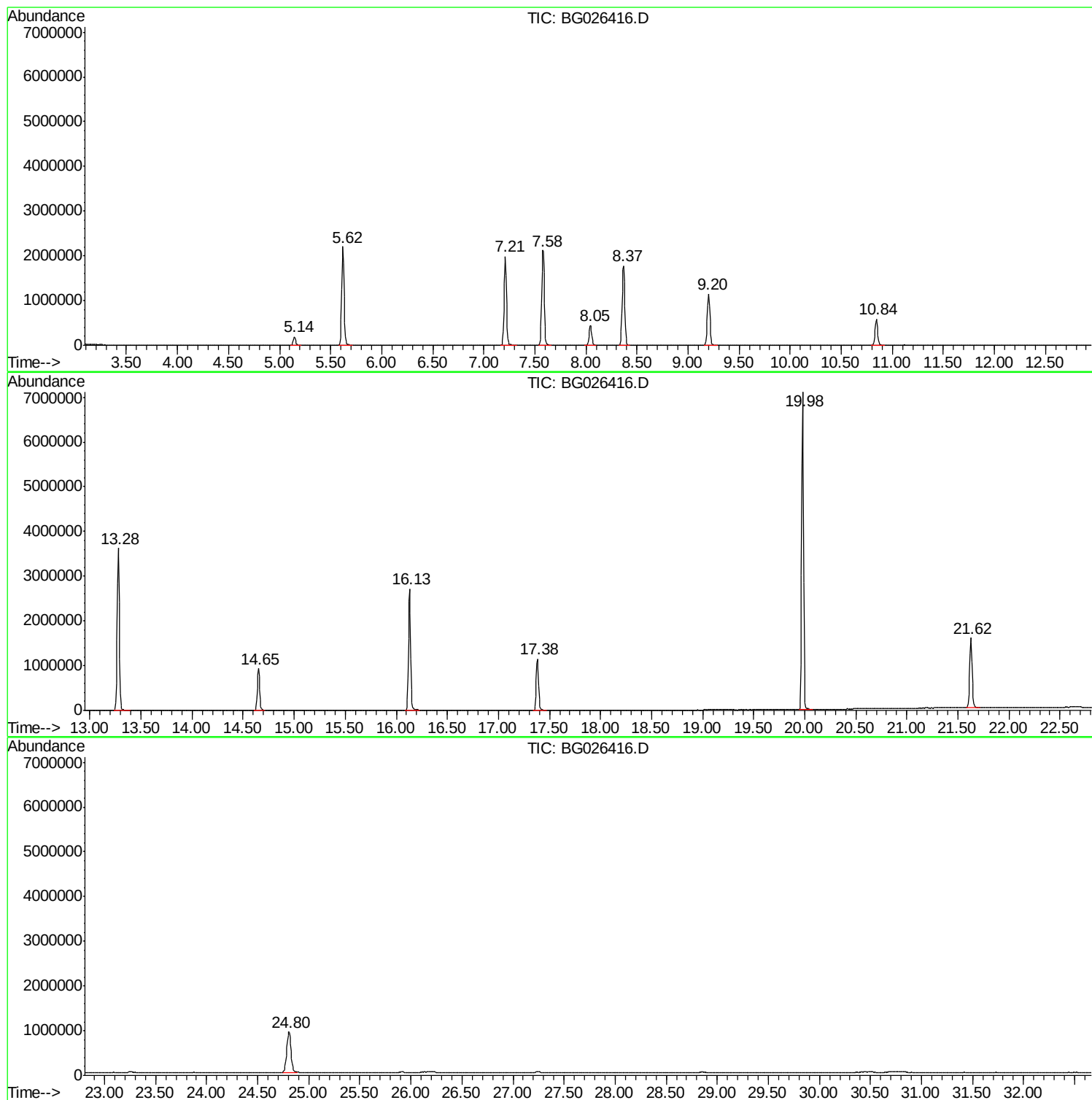
Sum of corrected areas: 45820297

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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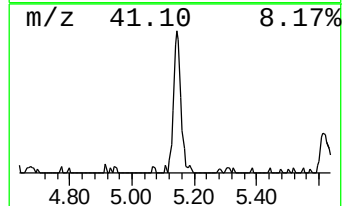
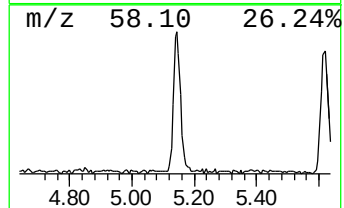
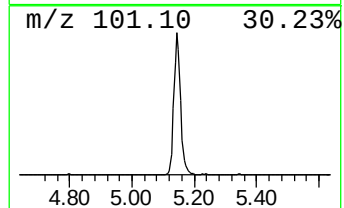
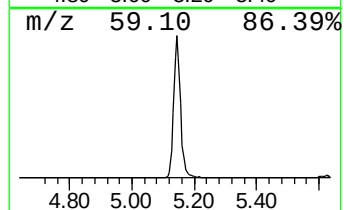
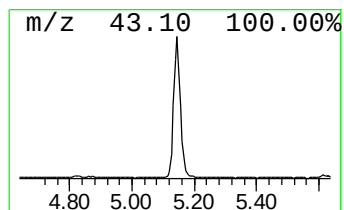
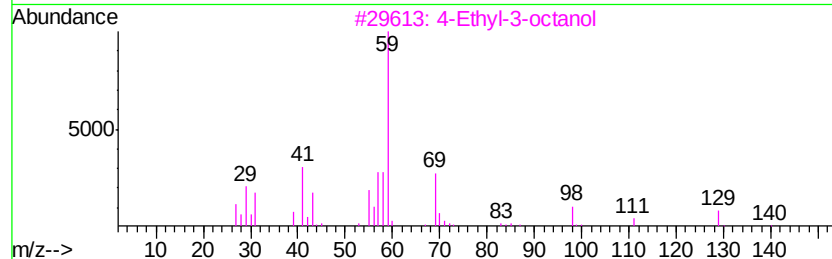
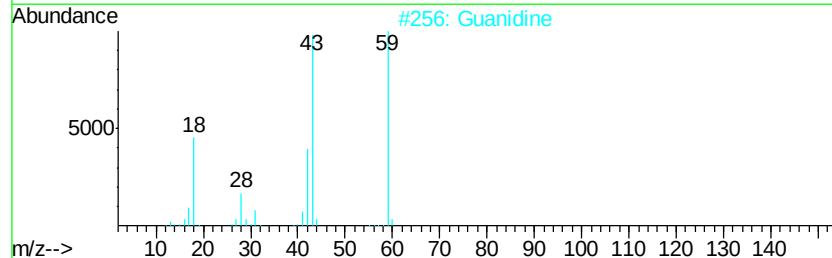
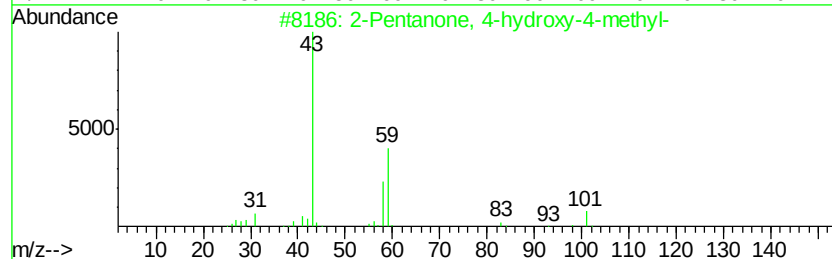
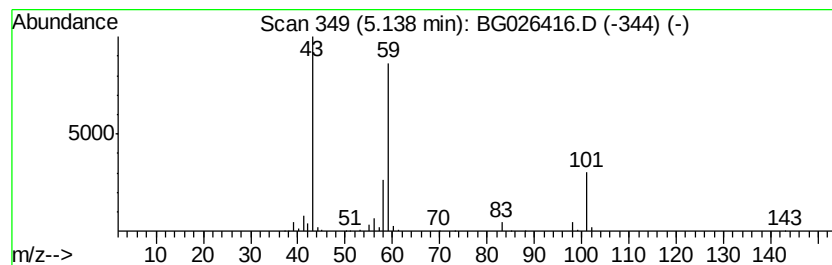
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	7.80 ng	296678	1,4-Dichlorobenzene-d4	8.04

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2			Guanidine	59	CH5N3	000113-00-8	47
3			4-Ethyl-3-octanol	158	C10H22O	019781-26-1	33
4			Succinic acid, heptyl 2-methoxye...	274	C14H26O5	1000325-80-7	33
5			1,8-Nonanediol, 8-methyl-	174	C10H22O2	054725-73-4	28



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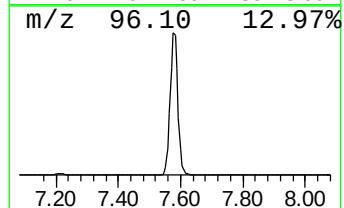
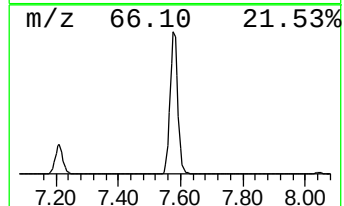
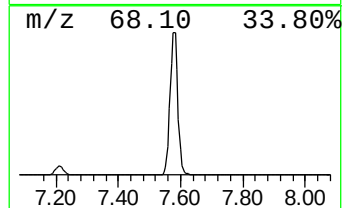
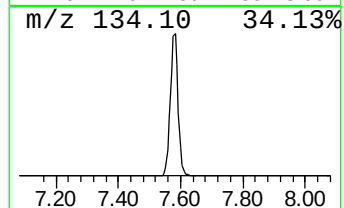
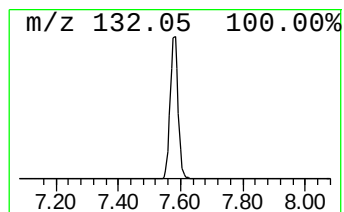
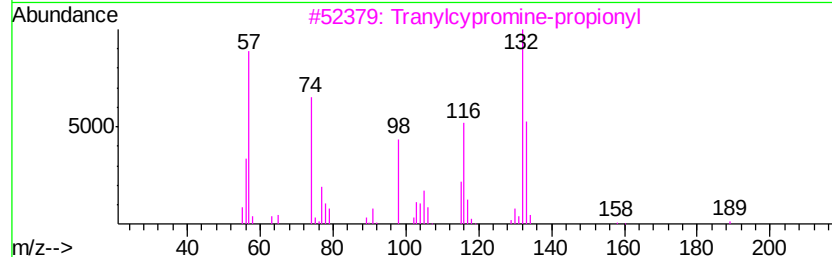
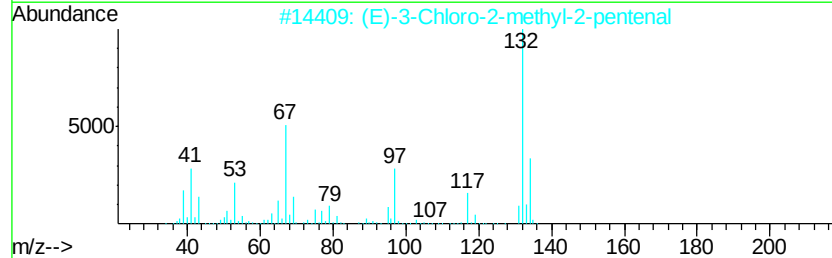
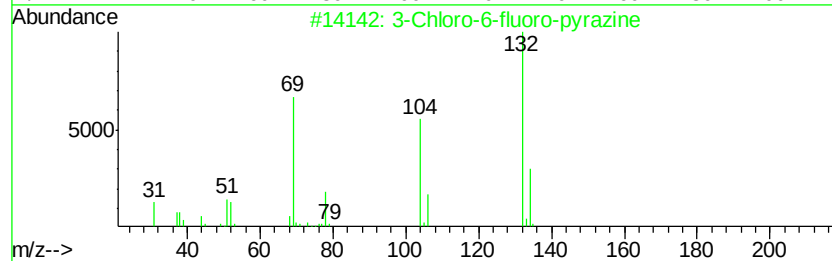
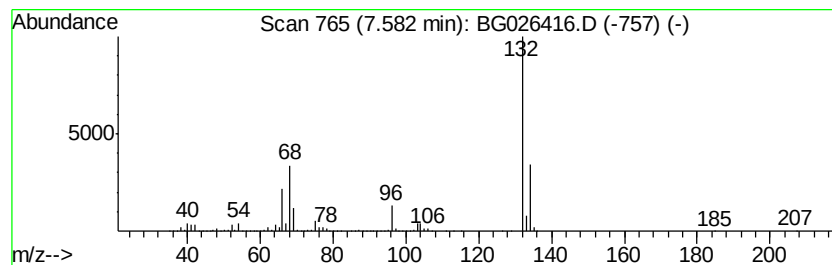
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Peak Number 2 unknown7.58 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.58	98.70 ng	3753550	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	38
2		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	17
3		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
4		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9
5		1H-Benzimidazole, 2-methyl-	132	C8H8N2	000615-15-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	7.8	ng	296678	1	8.04	760579	20.0
unknown7.58	7.58	98.7	ng	3753550	1	8.04	760579	20.0