

Data Path : Z:\HPCHEM1\BNA G\DATA\BG072716\
 Data File : BG023279.D
 Acq On : 27 Jul 2016 12:23
 Operator : UM/SJ
 Sample : PB92399BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB92399BL

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-BG072616.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	5.200	397	402	410	rBV	522970	798495	7.83%	1.544%
2	5.681	477	484	495	rBV	1713968	2782712	27.28%	5.381%
3	7.297	752	759	772	rBV	1710335	2974678	29.16%	5.752%
4	7.673	814	823	834	rBV	1668608	2904756	28.48%	5.617%
5	8.149	896	904	913	rVB	661183	1159657	11.37%	2.242%
6	8.477	952	960	969	rBV	1871118	3305476	32.41%	6.392%
7	9.323	1097	1104	1113	rBV	1495724	2728153	26.75%	5.275%
8	10.986	1380	1387	1399	rVB	951643	1668861	16.36%	3.227%
9	13.430	1795	1803	1814	rBV	4213696	6594848	64.66%	12.752%
10	14.816	2032	2039	2049	rVB2	1512241	2486094	24.37%	4.807%
11	16.308	2285	2293	2311	rBV	2097095	3239847	31.76%	6.265%
12	17.577	2502	2509	2520	rBV2	2140777	3366214	33.00%	6.509%
13	20.185	2946	2953	2959	rBV	7576879	10199957	100.00%	19.723%
14	21.895	3236	3244	3253	rBV2	2243493	3829001	37.54%	7.404%
15	25.308	3812	3825	3839	rBV3	1211331	3677972	36.06%	7.112%

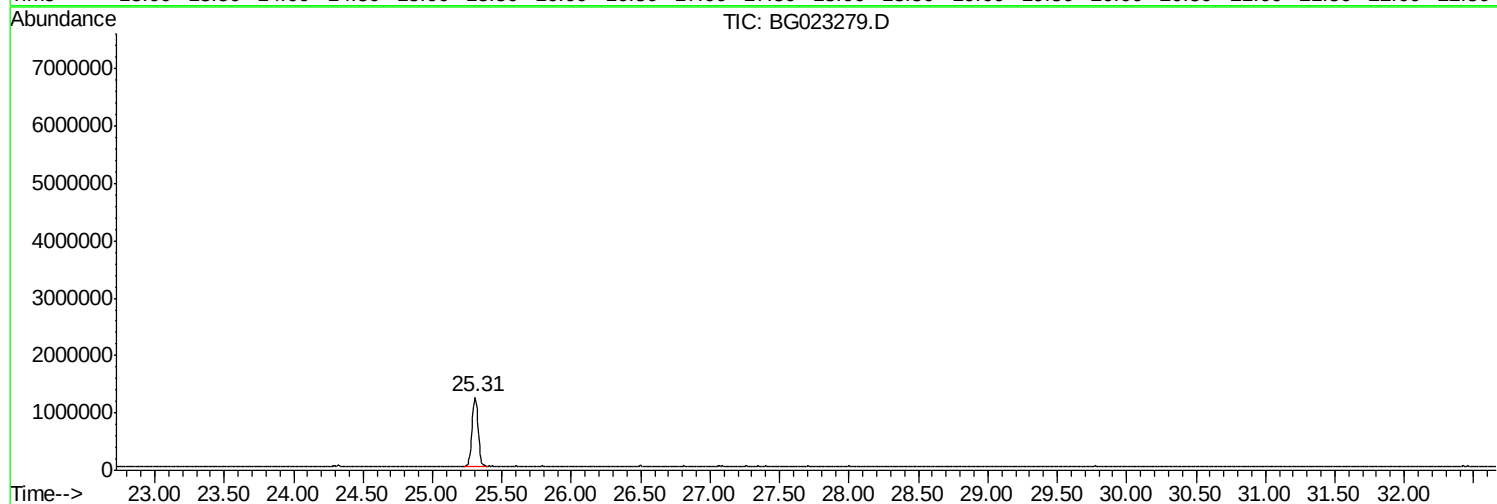
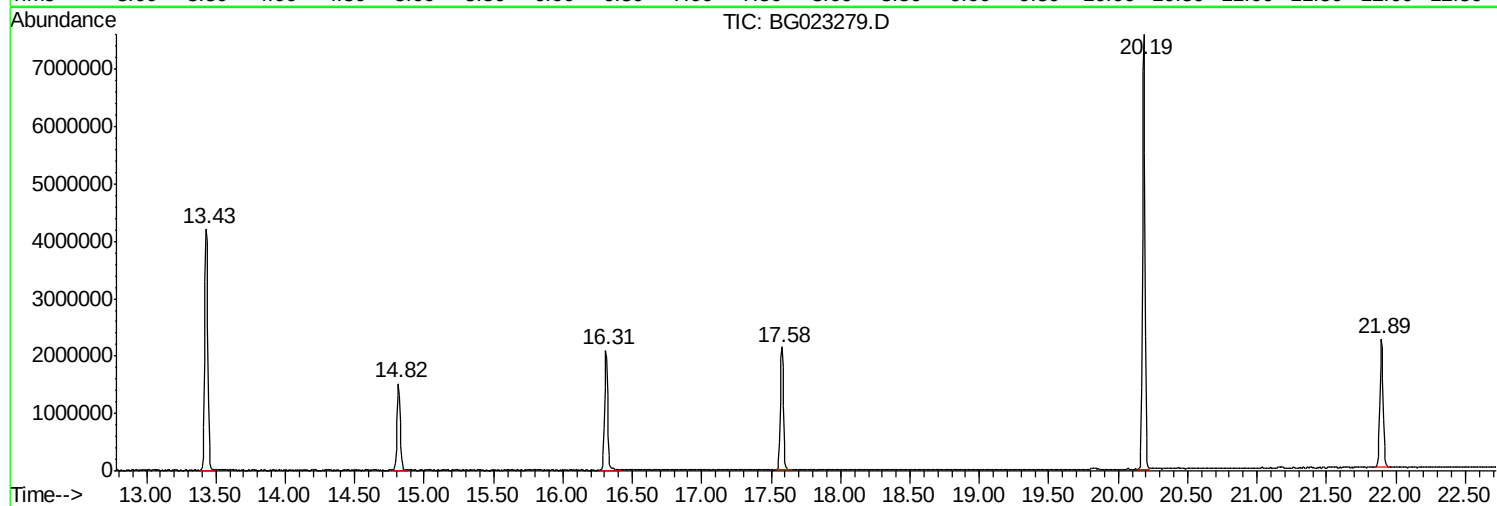
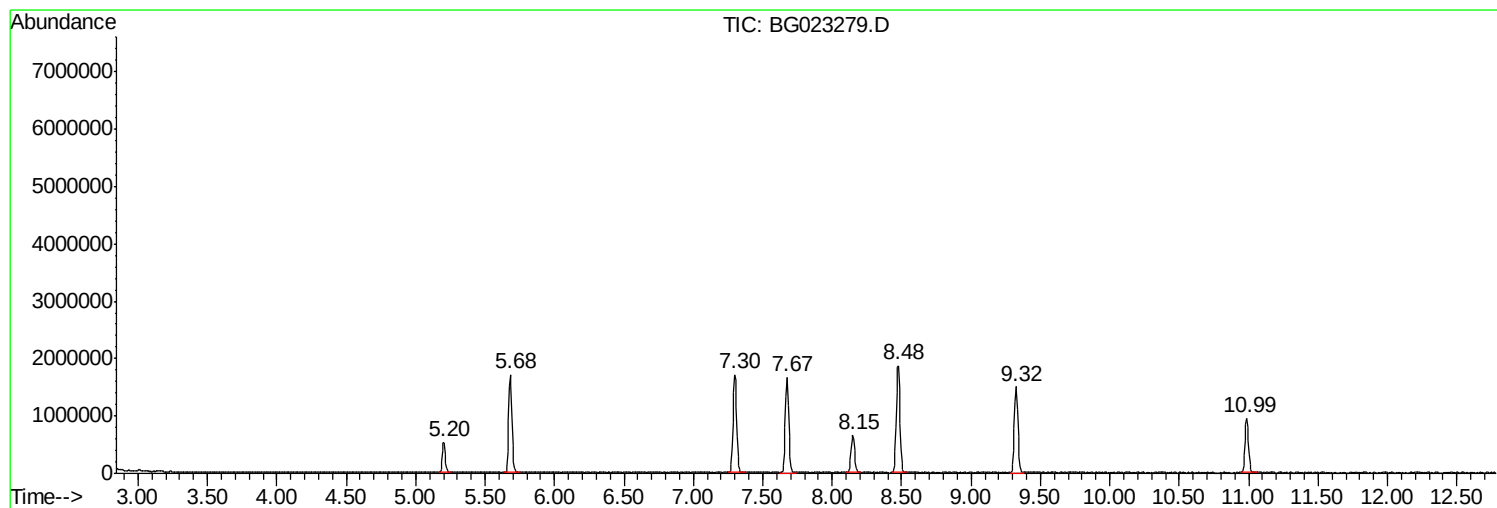
Sum of corrected areas: 51716721

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG072616.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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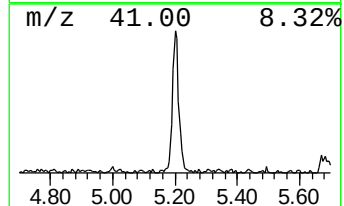
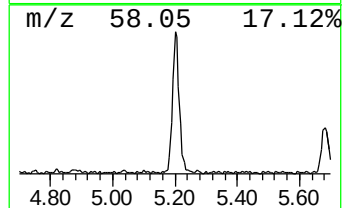
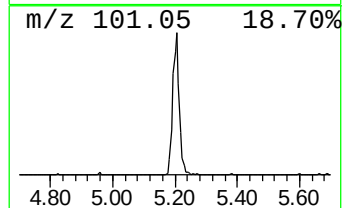
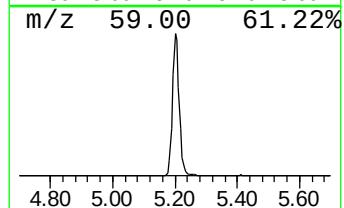
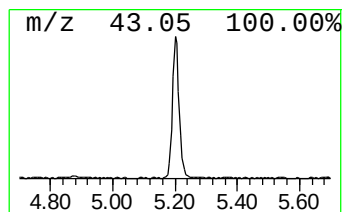
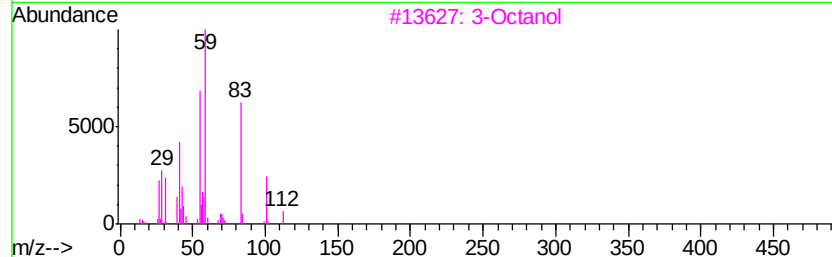
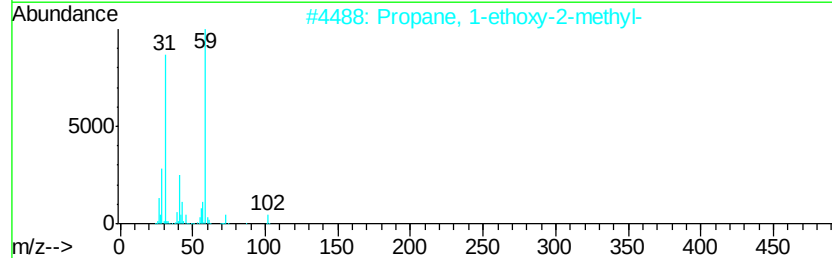
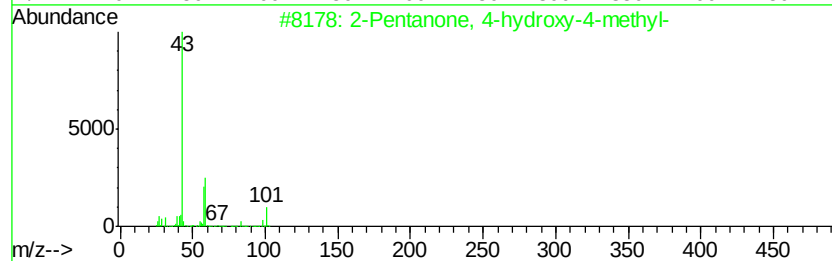
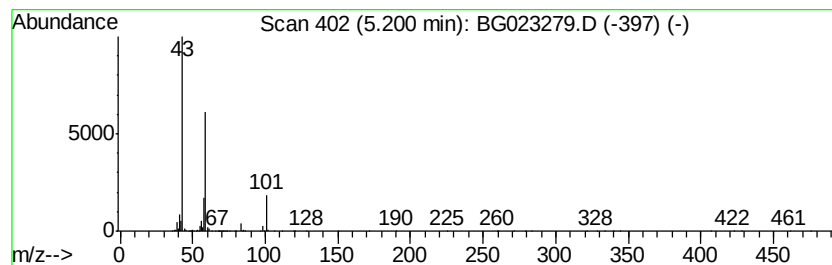
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.20	13.77 ng	798495	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64
2		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	32
3		3-Octanol	130	C8H18O	000589-98-0	28
4		1-Propanol, 2-(2-hydroxypropoxy)-	134	C6H14O3	000106-62-7	25
5		Acetic acid, cyano-, 1,1-dimethyl-	141	C7H11NO2	001116-98-9	23



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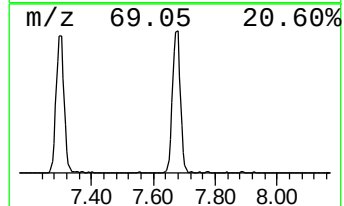
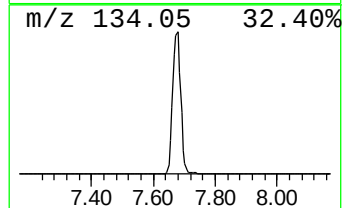
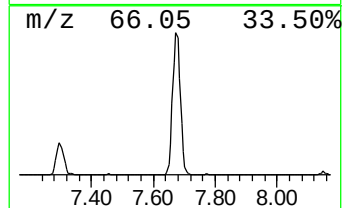
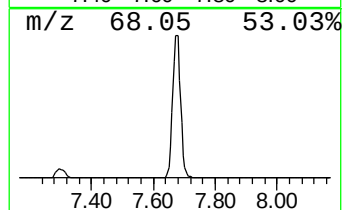
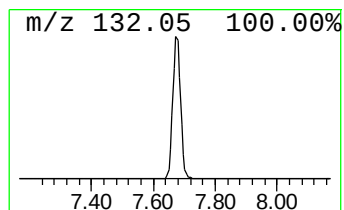
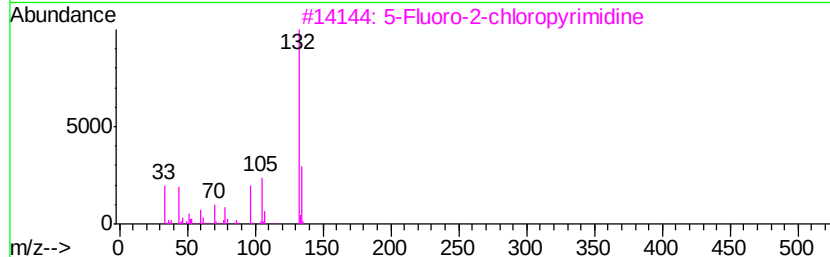
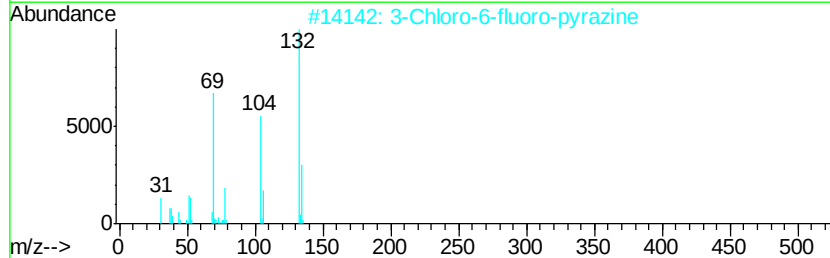
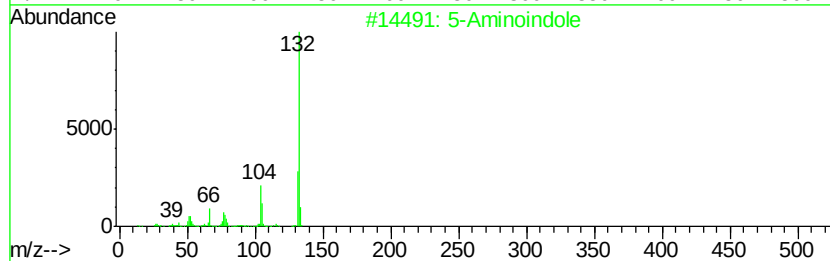
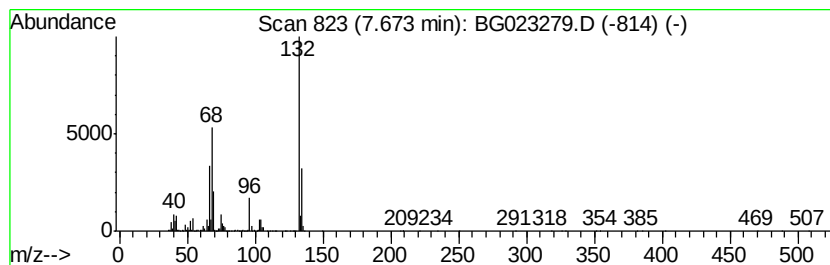
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Peak Number 2 unknown7.67 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.67	50.10 ng	2904760	1,4-Dichlorobenzene-d4	8.15

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Aminoindole	132	C8H8N2	005192-03-0	22
2		3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	16
3		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
4		2-Cyclohexen-1-one	96	C6H8O	000930-68-7	11
5		Tranlylcypromine-propionyl	189	C12H15NO	1000123-86-3	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.20	13.8	ng	798495	1	8.15	1159660	20.0
unknown7.67	7.67	50.1	ng	2904760	1	8.15	1159660	20.0