

Data Path : Z:\HPCHEM1\BNA G\Data\BG032317\
 Data File : BG026401.D
 Acq On : 23 Mar 2017 13:59
 Operator : SJ/MA
 Sample : PB97769BL
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB97769BL

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	5.137	343	349	361	rVB	250337	377383	4.91%	1.001%
2	5.613	423	430	441	rBV	1621966	2633647	34.26%	6.989%
3	7.200	692	700	713	rBV	1492183	2530192	32.92%	6.714%
4	7.570	756	763	774	rBV	1637789	2843596	36.99%	7.546%
5	8.040	836	843	852	rBV	424295	721409	9.39%	1.914%
6	8.363	890	898	910	rBV	1365399	2339459	30.44%	6.208%
7	9.197	1032	1040	1053	rBV	868696	1542467	20.07%	4.093%
8	10.842	1310	1320	1332	rBV	541522	1016955	13.23%	2.699%
9	13.275	1727	1734	1749	rBV	2944752	4449050	57.88%	11.806%
10	14.650	1960	1968	1981	rBV2	839373	1377460	17.92%	3.655%
11	16.124	2212	2219	2237	rBV	2034415	3162917	41.15%	8.393%
12	17.382	2426	2433	2446	rBV2	1115541	1776418	23.11%	4.714%
13	19.979	2869	2875	2888	rBV	5724455	7686506	100.00%	20.398%
14	21.624	3148	3155	3168	rBV	1581334	2607416	33.92%	6.919%
15	24.802	3685	3696	3708	rBV2	915443	2618496	34.07%	6.949%

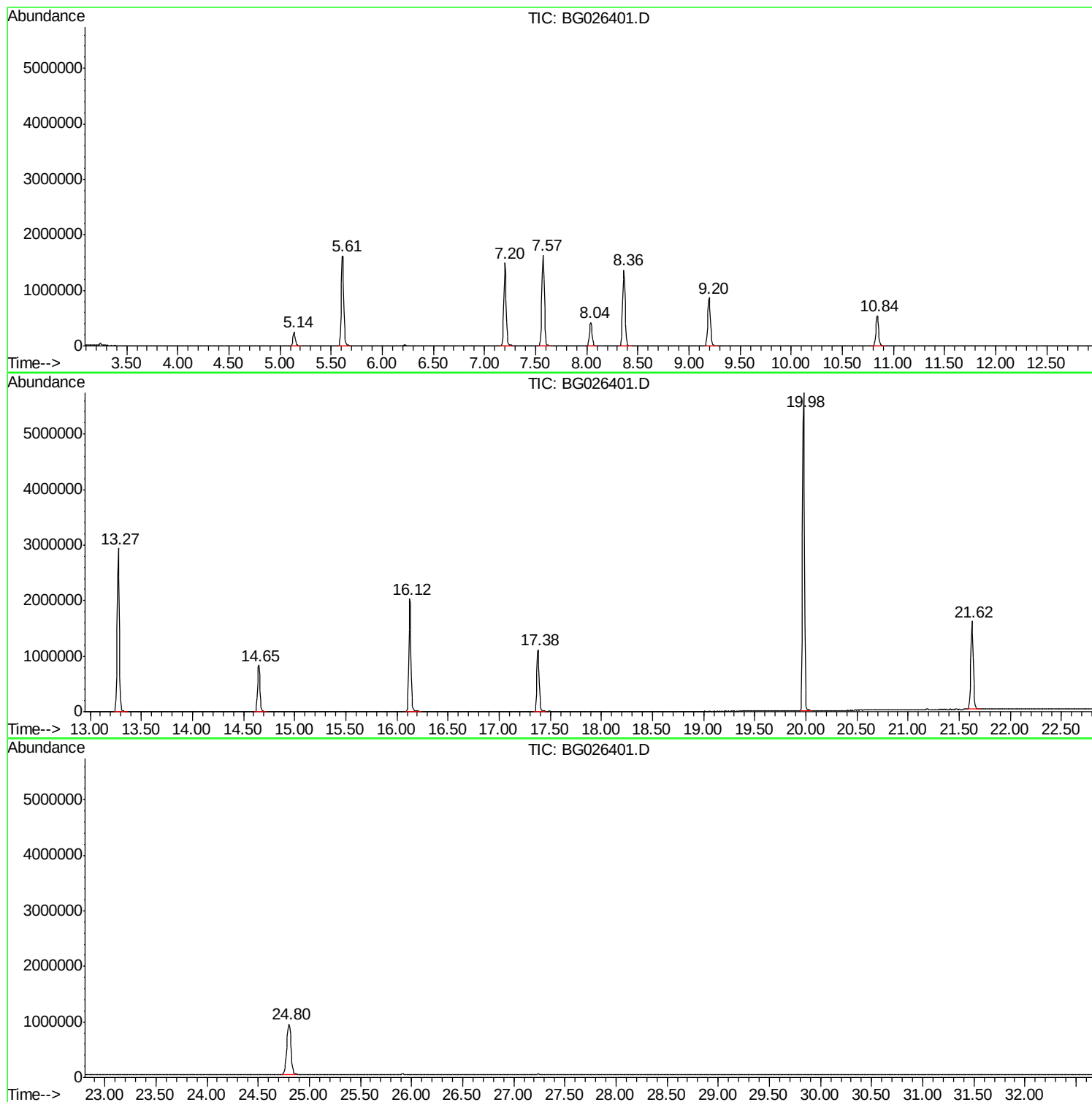
Sum of corrected areas: 37683371

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG032017.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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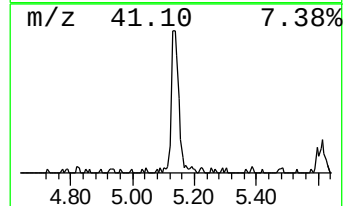
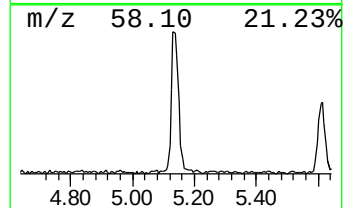
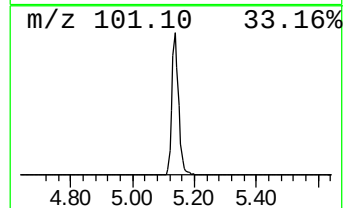
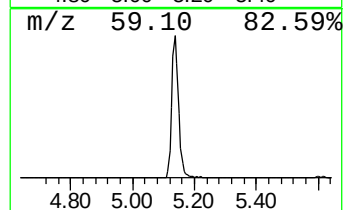
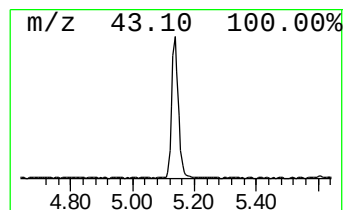
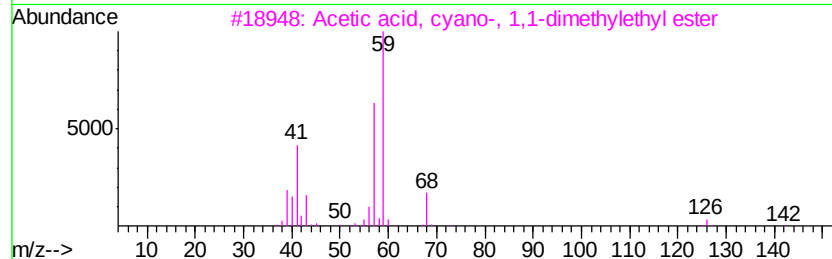
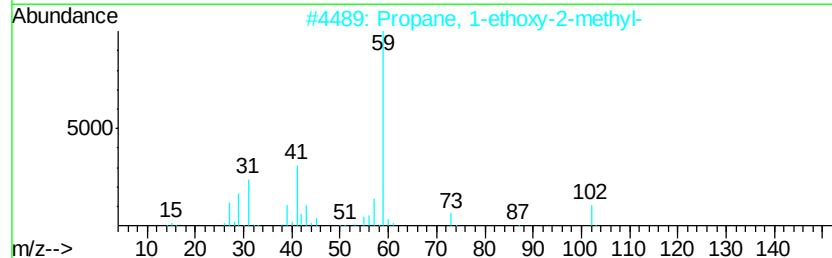
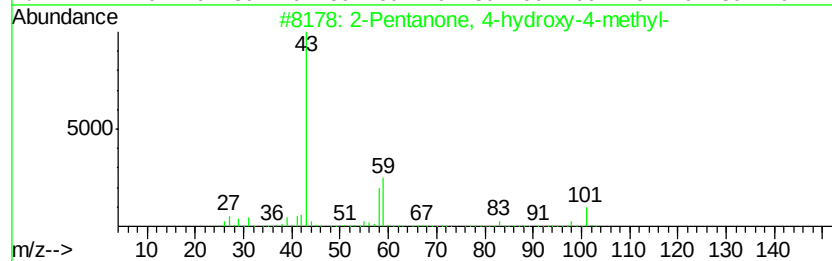
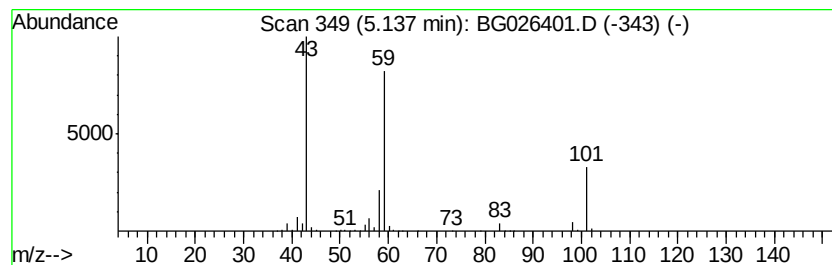
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Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.14	10.46 ng	377383	1,4-Dichlorobenzene-d4	8.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	25
3		Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
4		Guanidine	59	CH5N3	000113-00-8	9
5		5-Hexen-2-one	98	C6H10O	000109-49-9	9



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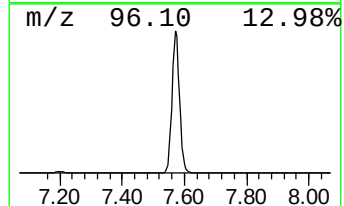
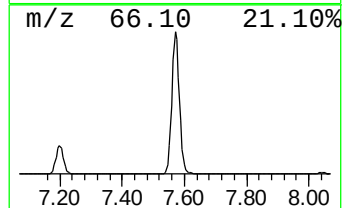
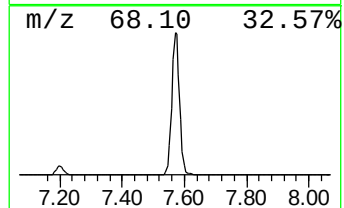
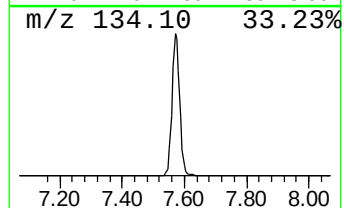
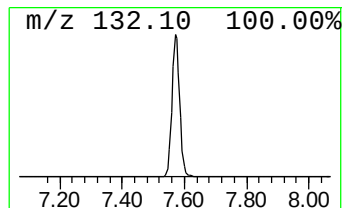
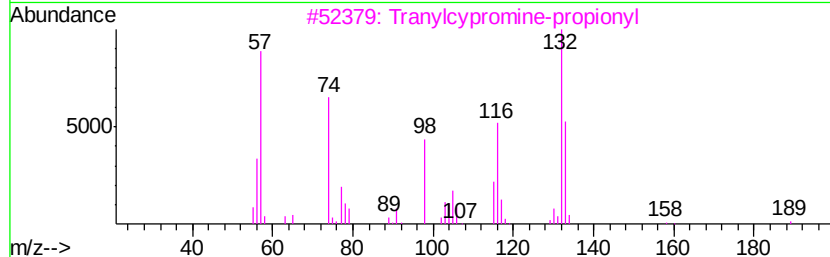
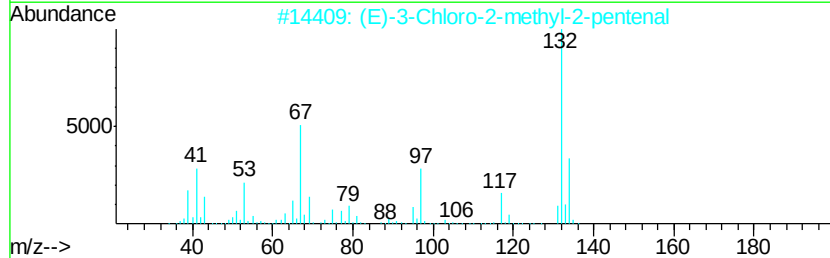
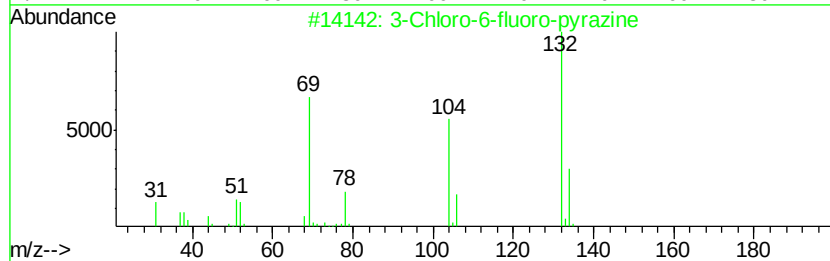
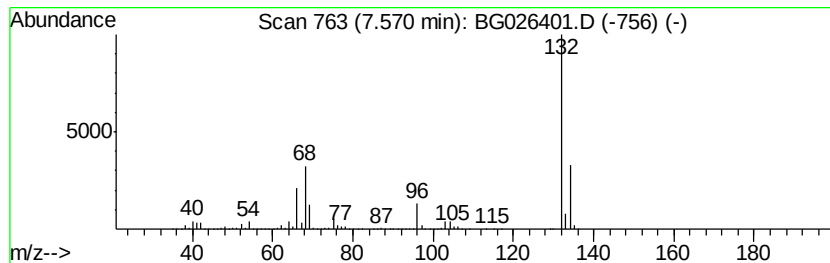
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Peak Number 2 unknown7.57 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.57	78.83 ng	2843600	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		Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	12
4		1H-Indazole, 7-methyl-	132	C8H8N2	1000316-00-7	10
5		4-Chloro-6-fluoro-pyrimidine	132	C4H2ClFN2	051422-01-6	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.14	10.5	ng	377383	1	8.04	721409	20.0
unknown7.57	7.57	78.8	ng	2843600	1	8.04	721409	20.0