

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060717\
 Data File : BG027298.D
 Acq On : 8 Jun 2017 00:32
 Operator : SJ/MA
 Sample : I3479-12
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SB-5(8-10)20170604

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-BG060517.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	5.700	336	342	351	rBV	69535	114869	4.06%	0.708%
2	6.141	410	417	433	rVB	687121	1158889	40.97%	7.144%
3	7.686	672	680	693	rBV	690670	1295384	45.80%	7.985%
4	8.092	741	749	758	rBV	657739	1200678	42.45%	7.401%
5	8.556	819	828	837	rBV	161821	298034	10.54%	1.837%
6	8.885	876	884	895	rVB	439482	811130	28.68%	5.000%
7	9.719	1016	1026	1038	rBV	365707	695229	24.58%	4.286%
8	11.376	1300	1308	1316	rBV	247570	476794	16.86%	2.939%
9	13.762	1707	1714	1723	rBV	1110815	1774800	62.75%	10.940%
10	14.572	1845	1852	1858	rBV	105404	164835	5.83%	1.016%
11	15.136	1940	1948	1959	rBV2	413164	710838	25.13%	4.382%
12	15.941	2080	2085	2091	rVB	22365	30573	1.08%	0.188%
13	16.617	2194	2200	2216	rVB	900997	1437385	50.82%	8.860%
14	16.805	2227	2232	2241	rBV	24747	44549	1.58%	0.275%
15	17.886	2409	2416	2424	rBV	580601	922030	32.60%	5.684%
16	18.726	2555	2559	2567	rBV3	33782	53126	1.88%	0.327%
17	20.465	2848	2855	2862	rBV	2107887	2828326	100.00%	17.435%
18	21.834	3083	3088	3095	rBV4	36872	67720	2.39%	0.417%
19	22.275	3155	3163	3171	rBV	568795	1100793	38.92%	6.786%
20	25.959	3779	3790	3806	rVB2	310086	1036482	36.65%	6.389%

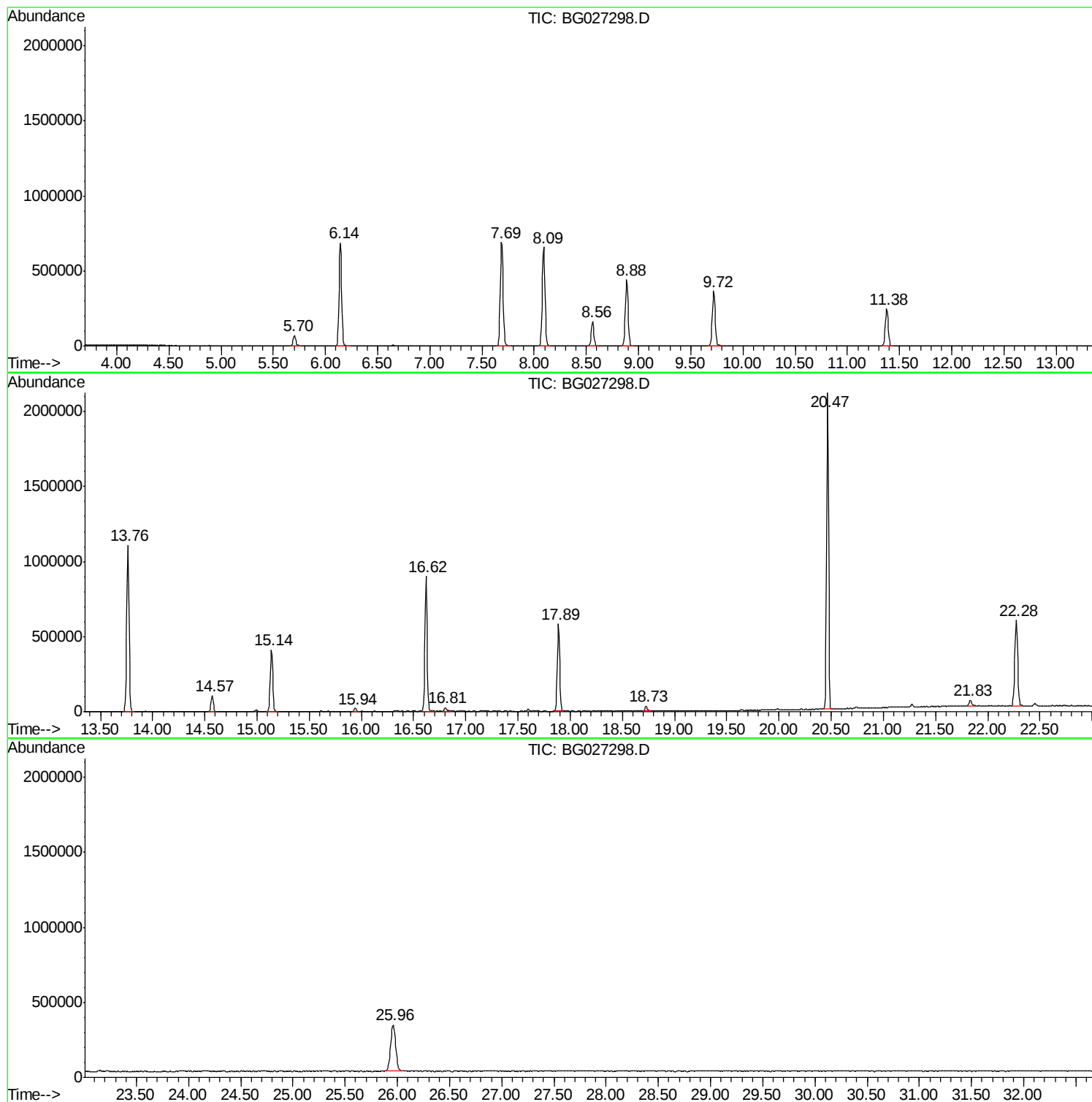
Sum of corrected areas: 16222464

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Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060517.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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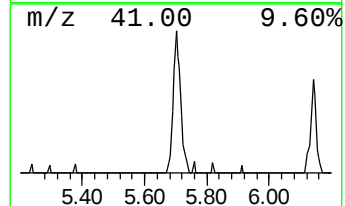
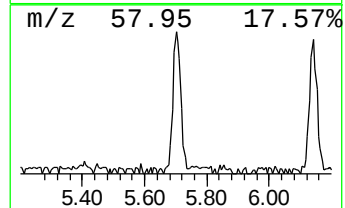
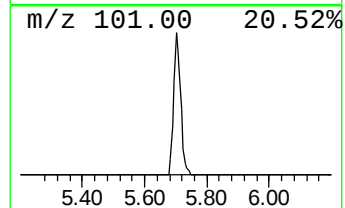
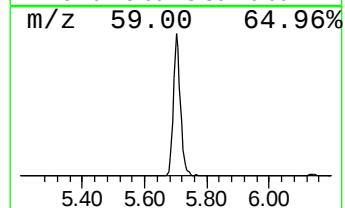
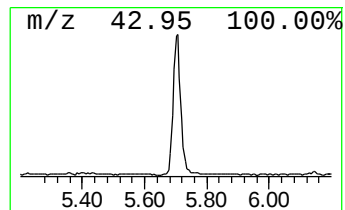
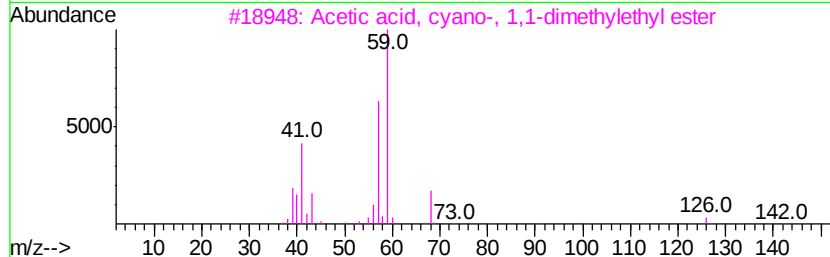
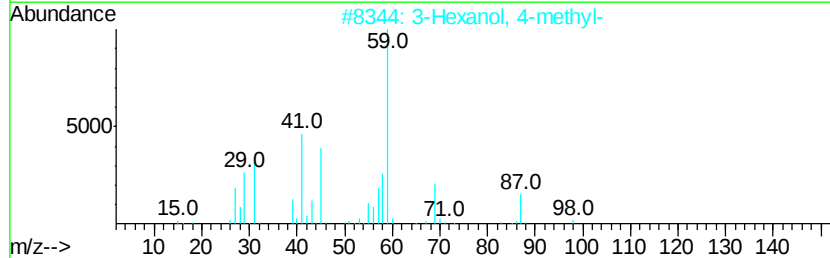
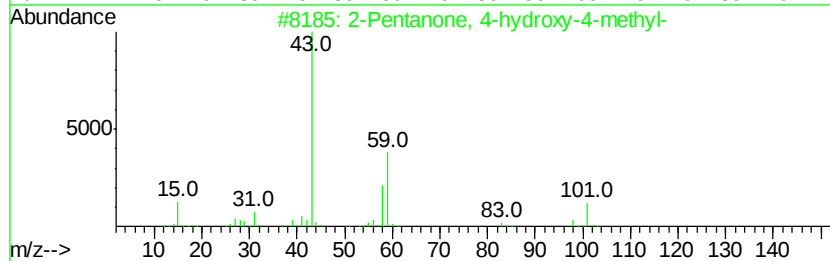
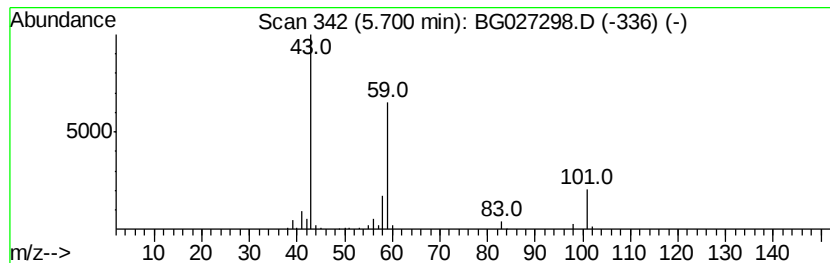
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TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.70	7.71 ng	114869	1,4-Dichlorobenzene-d4	8.56

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	25
3			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17
4			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	12
5			Acetone	58	C3H6O	000067-64-1	10



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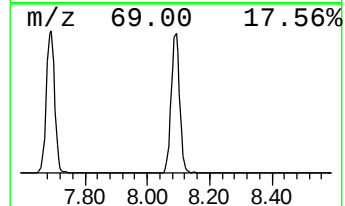
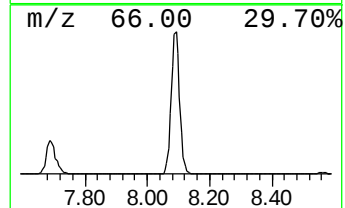
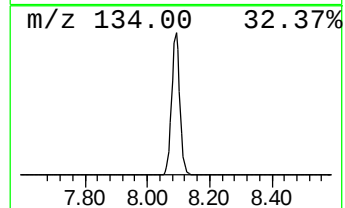
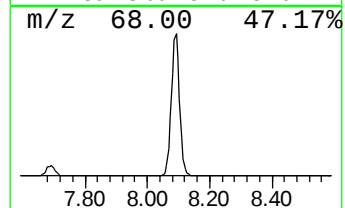
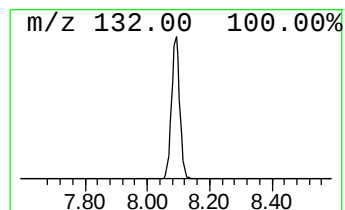
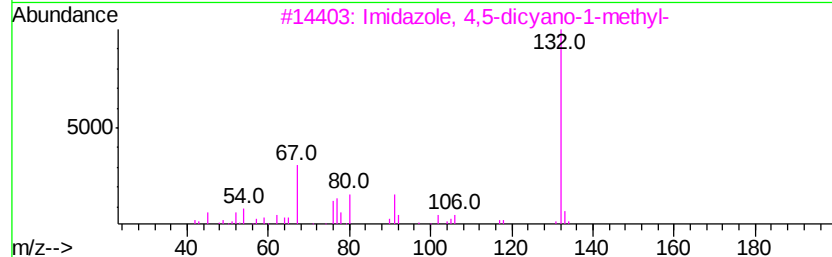
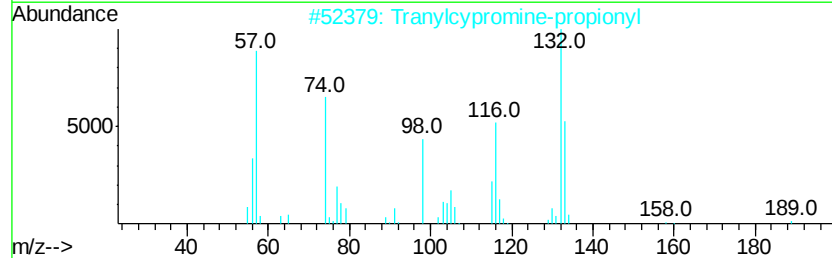
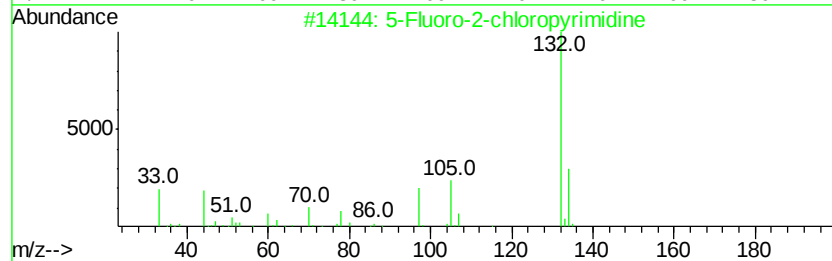
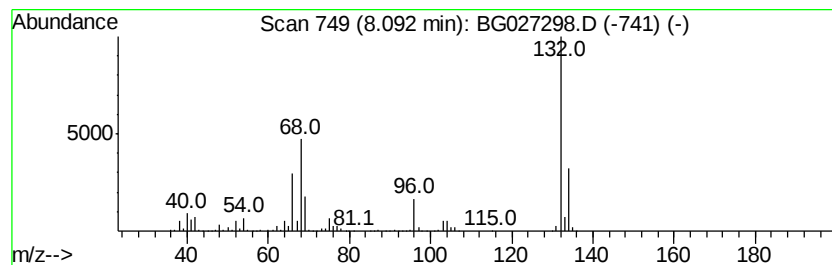
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Peak Number 2 unknown8.09 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.09	80.57 ng	1200680	1,4-Dichlorobenzene-d4	8.56

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	10
3		Imidazole, 4,5-dicyano-1-methyl-	132	C6H4N4	019485-35-9	9
4		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9
5		Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.70	7.7	ng	114869	1	8.56	298034	20.0
unknown8.09	8.09	80.6	ng	1200680	1	8.56	298034	20.0