

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121415\
 Data File : BF083797.D
 Acq On : 15 Dec 2015 4:29
 Operator : UM/IZ
 Sample : G4779-12
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 W-51

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_F\METHODS\8270-BF121315.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	5.070	259	261	264	rBV	157334	170807	4.69%	0.760%
2	5.493	296	298	301	rBV	914618	1004161	27.58%	4.467%
3	6.521	385	388	390	rBV	830731	733992	20.16%	3.265%
4	6.659	398	400	403	rBV	1440095	1933985	53.11%	8.603%
5	6.899	419	421	423	rBB	759156	633909	17.41%	2.820%
6	7.059	432	435	437	rVB	1991231	1930888	53.03%	8.589%
7	7.459	467	470	472	rBV	1670977	1577820	43.33%	7.019%
8	8.179	531	533	535	rBV	816494	807068	22.16%	3.590%
9	9.265	625	628	630	rBV	3245490	3433470	94.29%	15.273%
10	9.939	684	687	689	rBV	1082740	1018549	27.97%	4.531%
11	10.373	724	725	727	rVB	42247	38926	1.07%	0.173%
12	10.590	741	744	748	rBV	14568	37138	1.02%	0.165%
13	10.728	753	756	758	rBV	1885058	2285678	62.77%	10.168%
14	10.853	765	767	773	rVB	36083	70512	1.94%	0.314%
15	11.299	802	806	807	rBV	39725	47023	1.29%	0.209%
16	11.413	814	816	819	rVB	987879	999327	27.44%	4.445%
17	11.631	832	835	837	rBV	104513	93894	2.58%	0.418%
18	12.785	934	936	939	rBV	109293	130300	3.58%	0.580%
19	13.002	952	955	958	rBV	3131826	3641247	100.00%	16.198%
20	13.779	1021	1023	1028	rBV2	54497	69490	1.91%	0.309%
21	13.905	1031	1034	1039	rBV3	28335	58569	1.61%	0.261%
22	14.042	1044	1046	1049	rBV	874599	1009749	27.73%	4.492%
23	15.482	1169	1172	1175	rBV	640695	753692	20.70%	3.353%

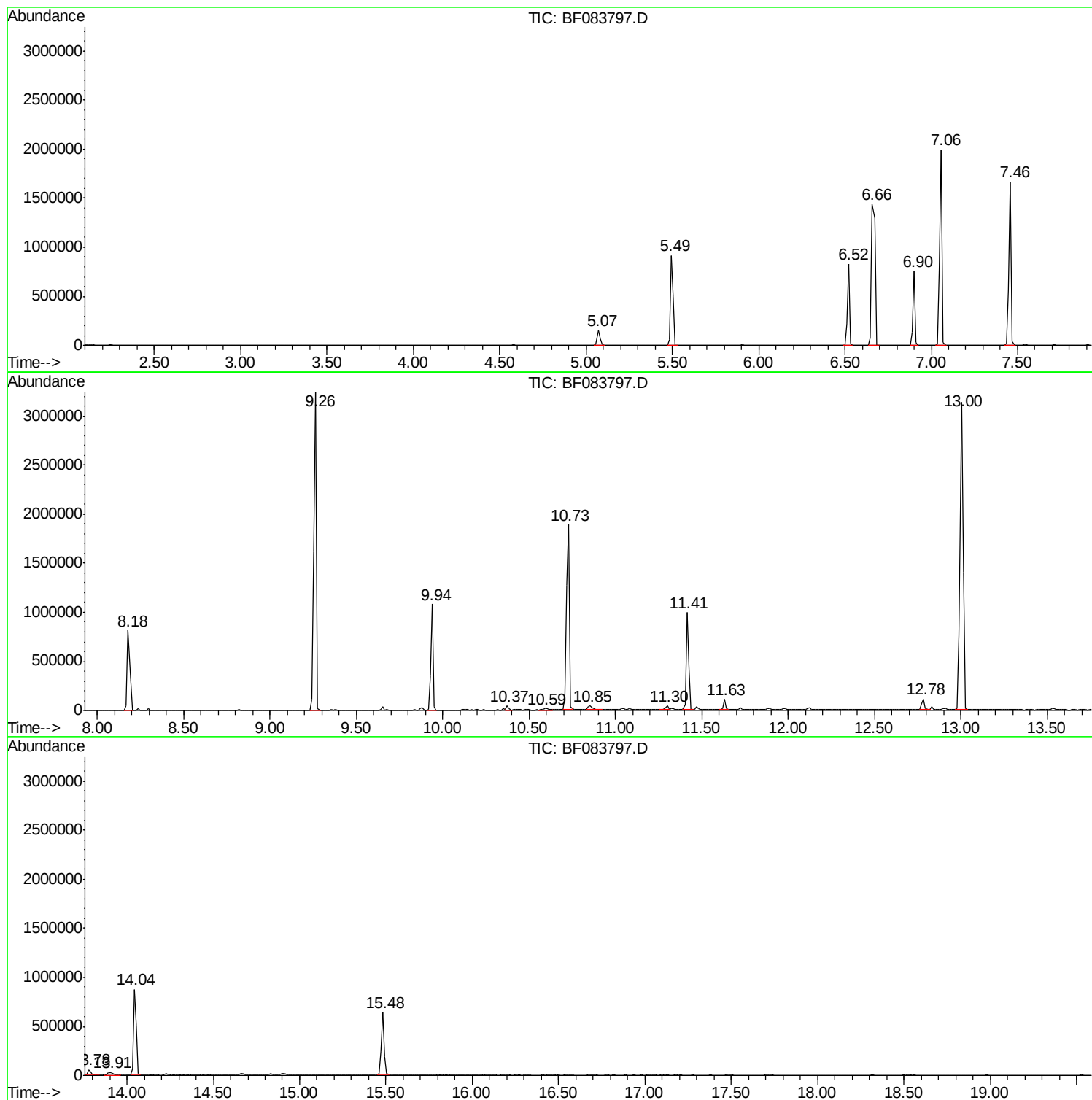
Sum of corrected areas: 22480194

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P



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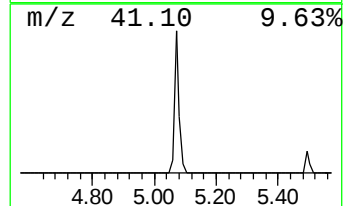
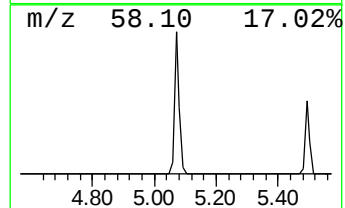
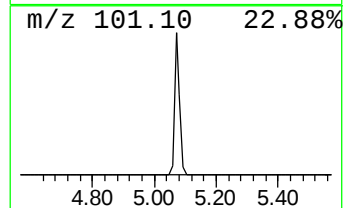
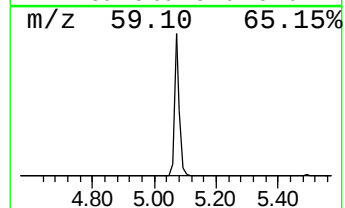
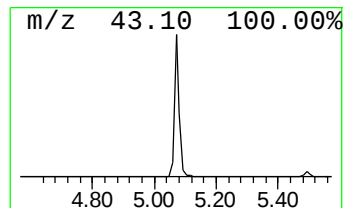
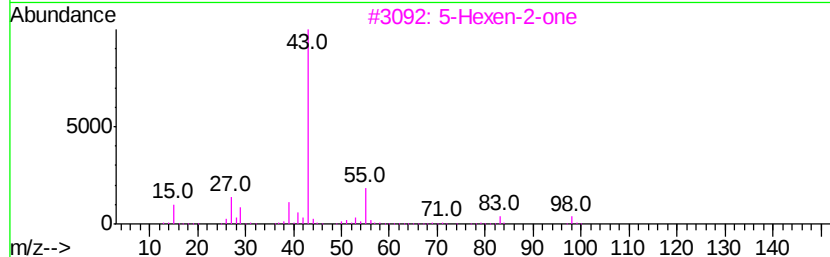
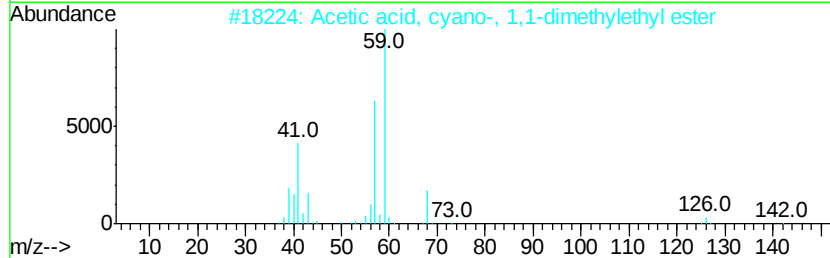
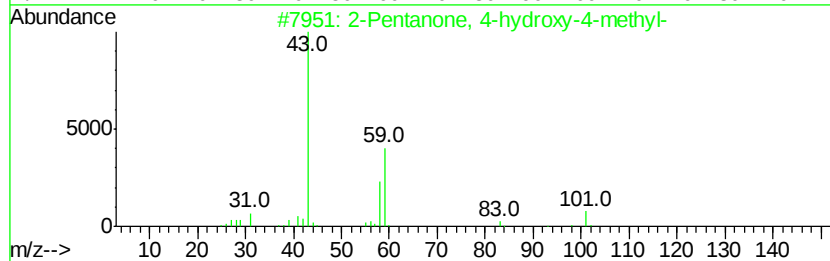
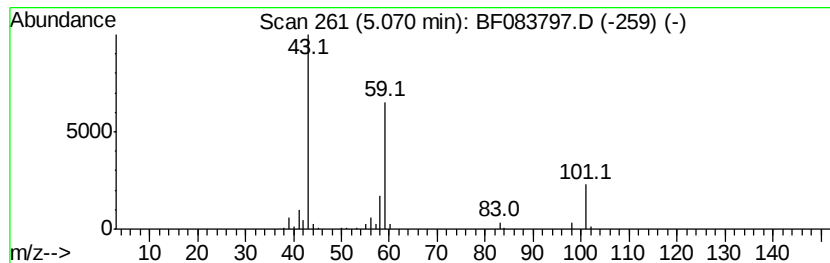
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
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.
5.07	5.39 ng	170807	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	25
3			5-Hexen-2-one	98	C6H10O	000109-49-9	9
4			Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9
5			Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9



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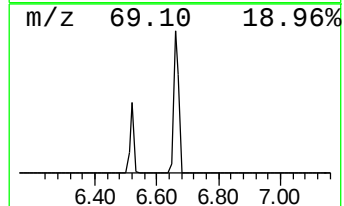
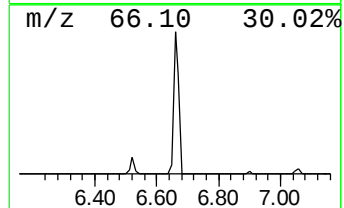
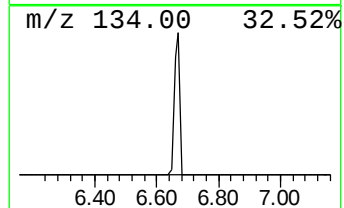
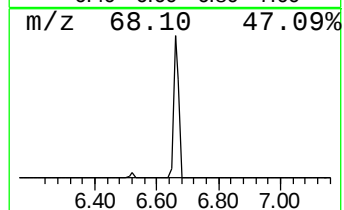
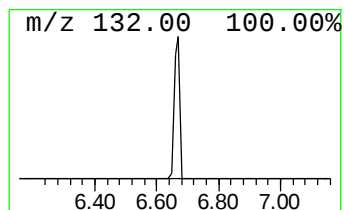
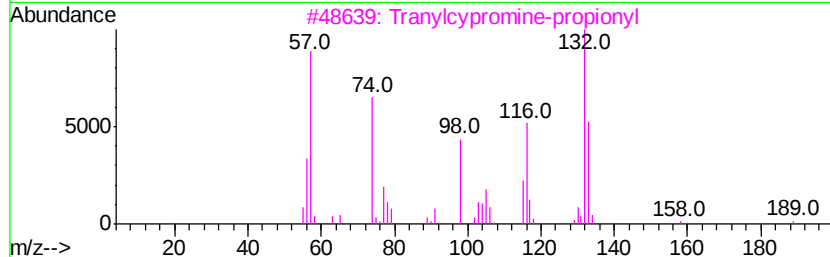
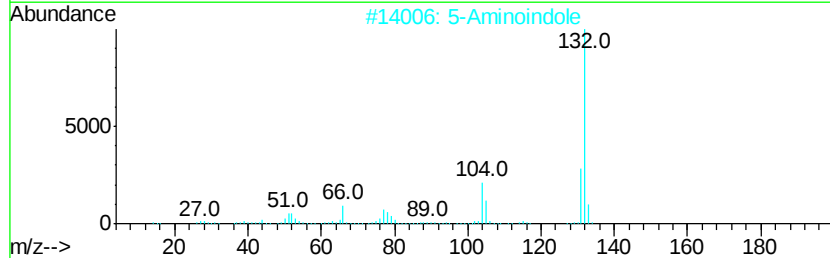
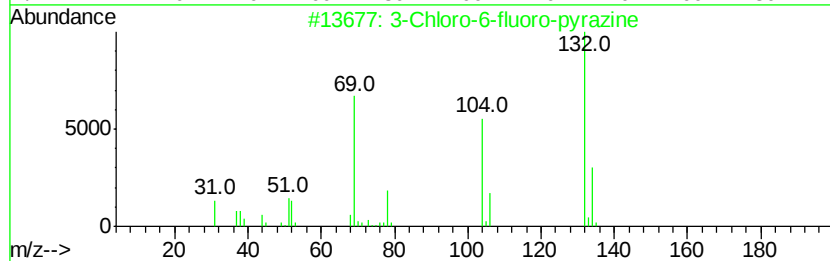
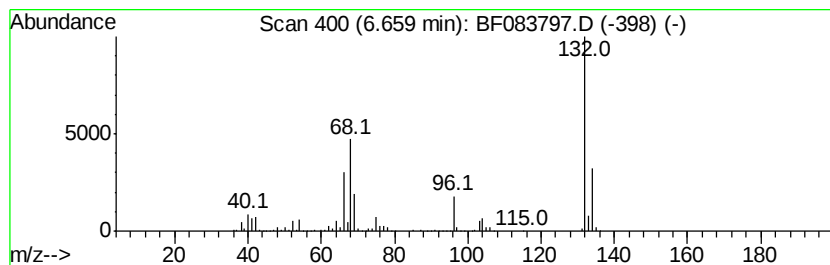
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Peak Number 2 unknown6.66 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.66	61.02 ng	1933990	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2			5-Aminoindole	132	C8H8N2	005192-03-0	12
3			Tranvlcypropine-propionyl	189	C12H15NO	1000123-86-3	10
4			Bicyclo[4.2.0]octa-1,3,5-triene,...	132	C10H12	028749-81-7	9
5			4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	9



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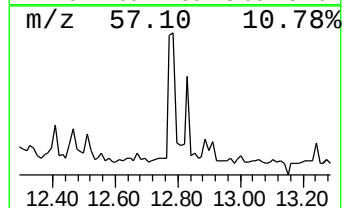
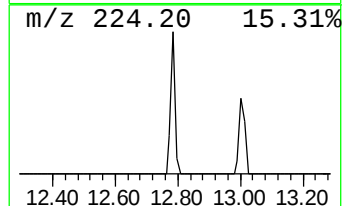
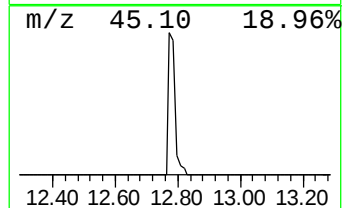
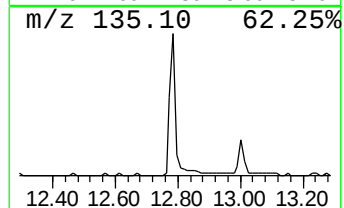
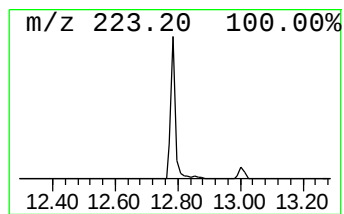
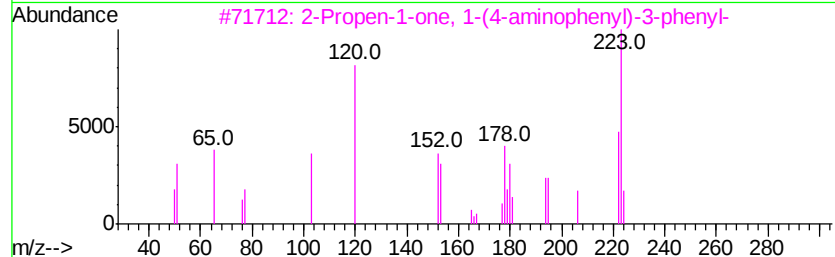
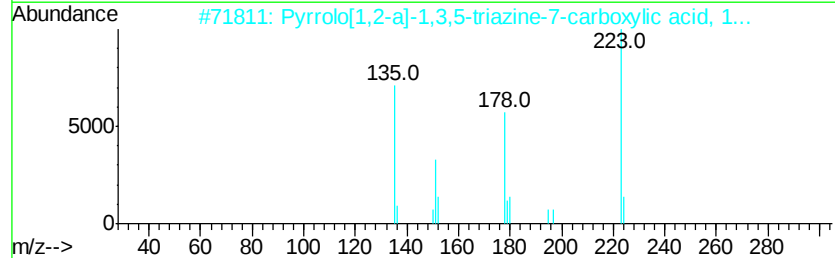
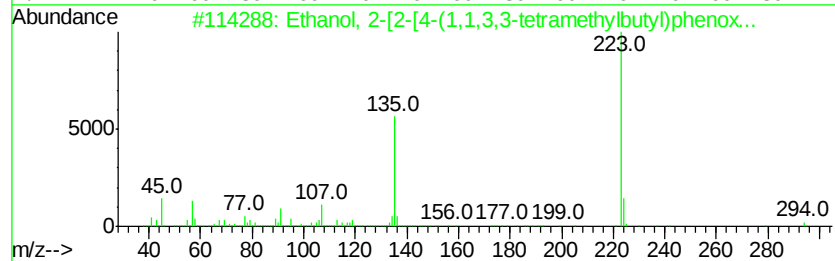
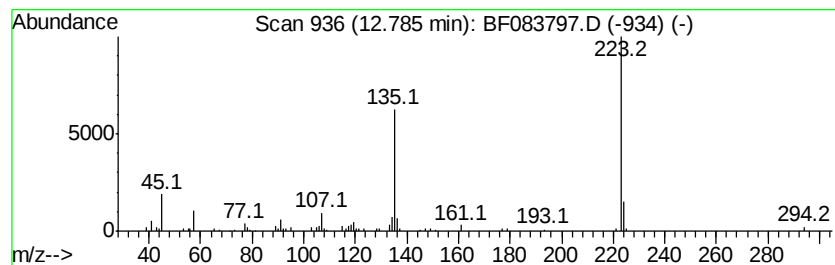
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Peak Number 4 Ethanol, 2-[2-[4-(1,1,3,3-t... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.79	2.58 ng	130300	Chrysene-d12	14.04

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Ethanol, 2-[2-[4-(1,1,3,3-tetram...	294	C18H30O3	002315-61-9	94
2		Pyrrolo[1,2-a]-1,3,5-triazine-7-...	223	C9H9N3O4	054449-89-7	90
3		2-Propen-1-one, 1-(4-aminophenyl)...	223	C15H13NO	002403-30-7	38
4		Benzenamine, N,N-dimethyl-4-(2-p...	223	C16H17N	001145-73-9	38
5		Benzenamine, N,N-dimethyl-4-(2-p...	223	C16H17N	000838-95-9	38



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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST02.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	5.07	5.4	ng	170807	1	6.90	633909	20.0
unknown6.66	6.66	61.0	ng	1933990	1	6.90	633909	20.0
Ethanol, 2-[2-[4-...	12.79	2.6	ng	130300	5	14.04	1009750	20.0