

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF080116\
 Data File : BF089519.D
 Acq On : 1 Aug 2016 19:26
 Operator : UM/SJ
 Sample : H4251-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-04

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-BF072716.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	1.655	4	6	58	rVB	893068	2351375	100.00%	22.340%
2	4.559	258	260	267	rVB	45962	76800	3.27%	0.730%
3	5.039	301	302	324	rBV9	316482	476763	20.28%	4.530%
4	6.136	396	398	405	rBV	212397	280236	11.92%	2.662%
5	6.239	405	407	422	rVB	715306	970813	41.29%	9.224%
6	6.479	426	428	434	rBV	179937	203209	8.64%	1.931%
7	6.627	439	441	444	rBV	767926	868925	36.95%	8.256%
8	7.050	476	478	497	rBV	613362	753526	32.05%	7.159%
9	7.759	538	540	552	rBV	221972	297573	12.66%	2.827%
10	8.845	632	635	637	rBV	1554192	1447410	61.56%	13.752%
11	9.508	691	693	695	rBV	349903	334377	14.22%	3.177%
12	10.296	759	762	764	rBV	605034	739488	31.45%	7.026%
13	10.982	819	822	824	rBV	217176	294551	12.53%	2.798%
14	12.559	957	960	963	rBV	1072061	1048844	44.61%	9.965%
15	13.599	1049	1051	1059	rVV	137383	192864	8.20%	1.832%
16	14.925	1165	1167	1177	rVB	154018	188592	8.02%	1.792%

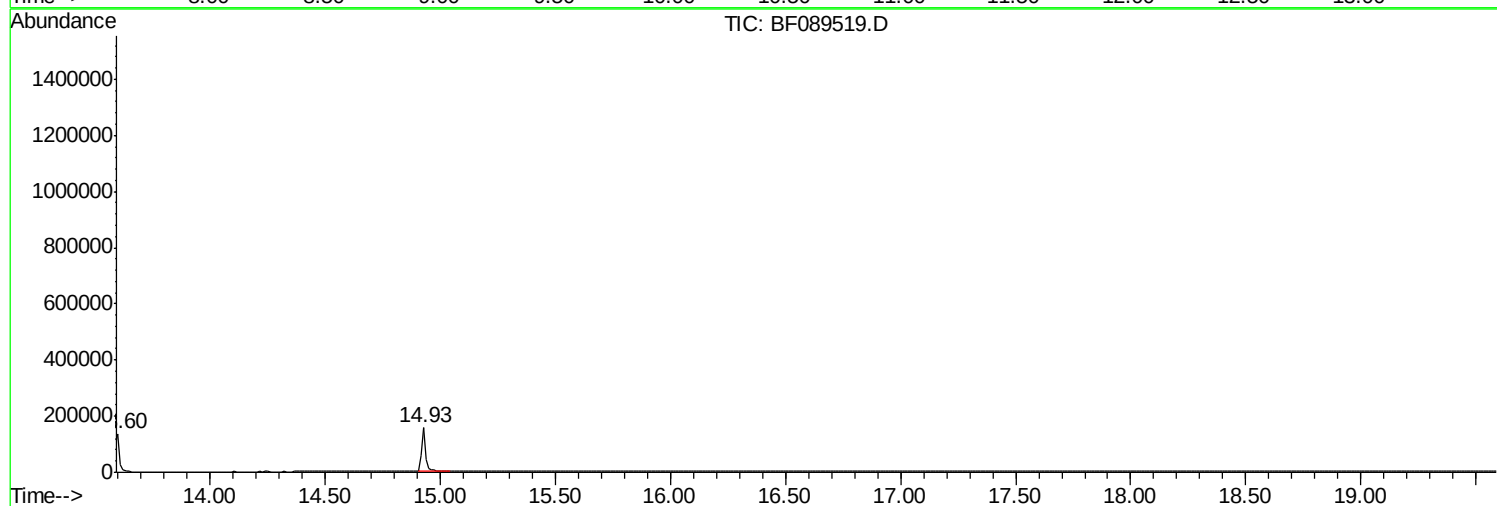
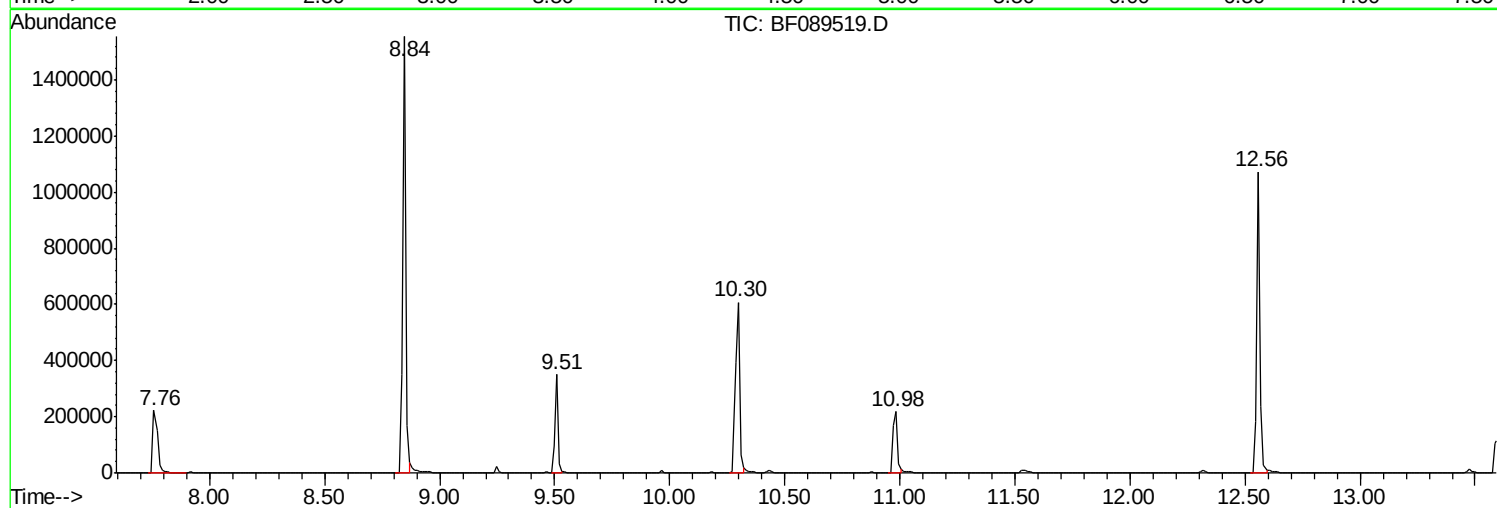
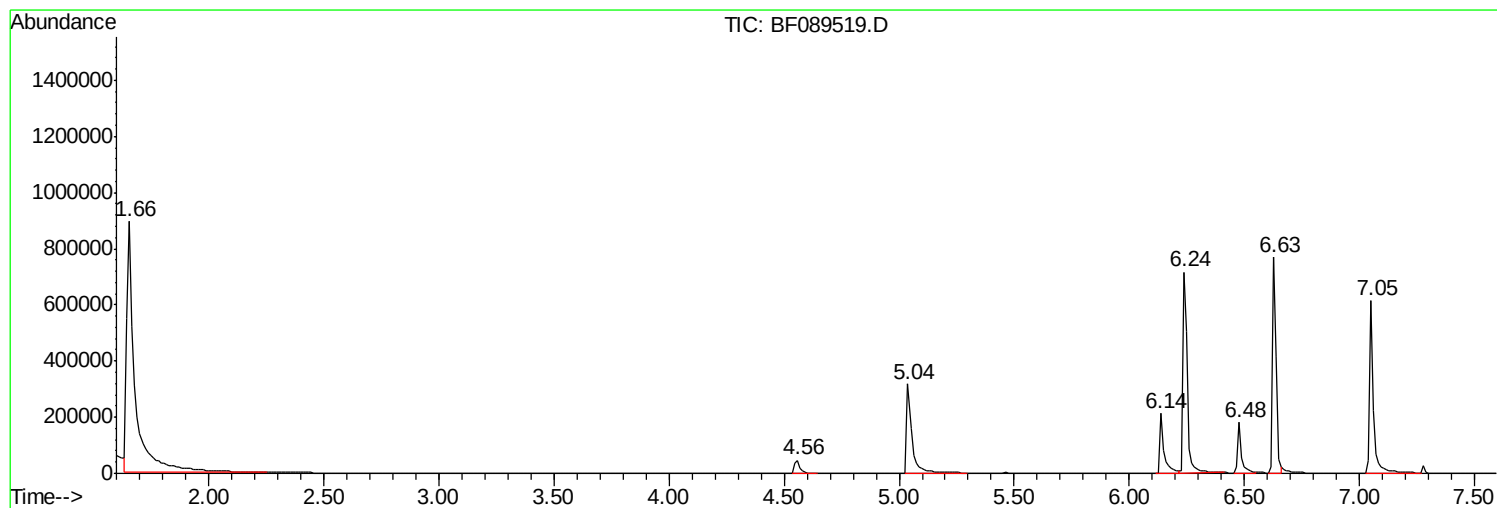
Sum of corrected areas: 10525346

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072716.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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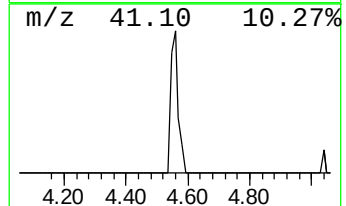
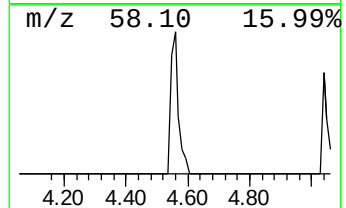
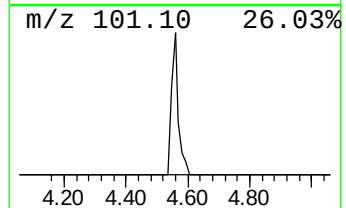
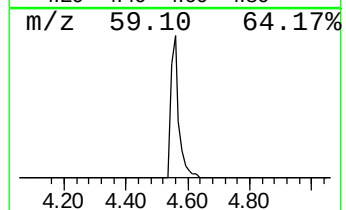
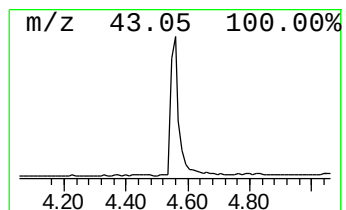
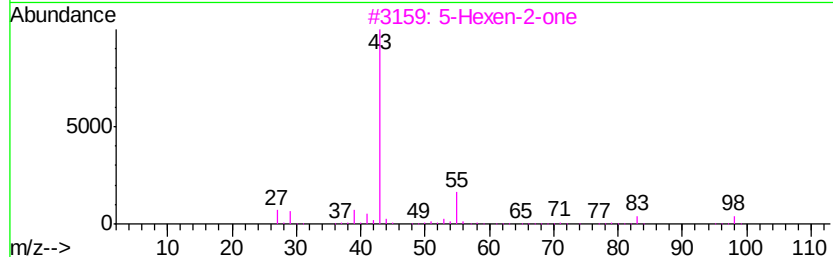
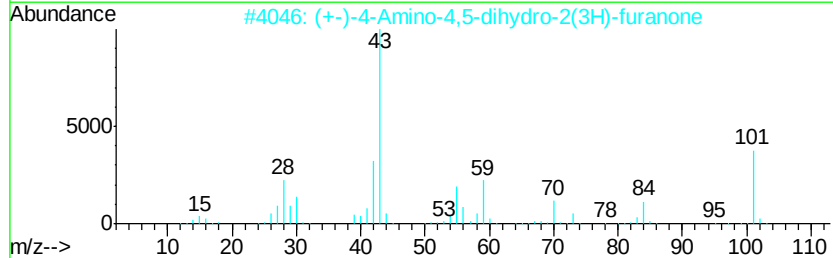
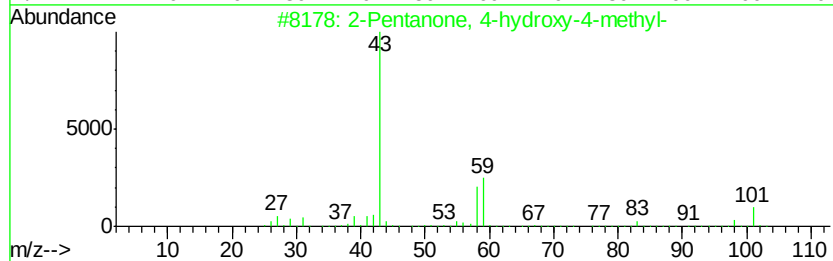
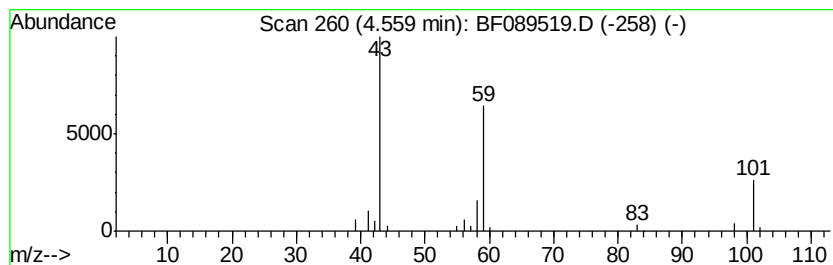
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Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.56	7.56 ng	76800	1,4-Dichlorobenzene-d4	6.48

Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	50
2		(+)-4-Amino-4,5-dihydro-2(3H)-furanone	101	C4H7NO2	016504-58-8	43
3		5-Hexen-2-one	98	C6H10O	000109-49-9	9
4		Propane, 1-ethoxy-2-methyl-	102	C6H14O	000627-02-1	9
5		Morpholine, 4-methyl-	101	C5H11NO	000109-02-4	9



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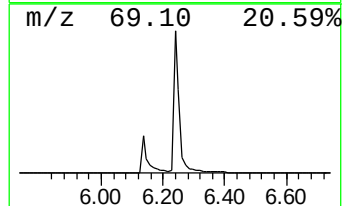
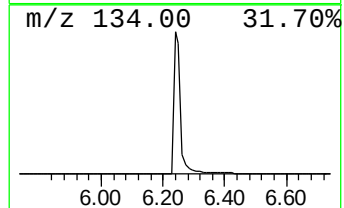
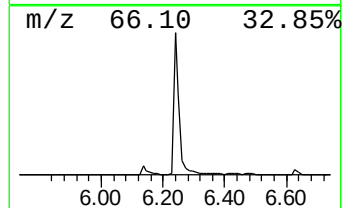
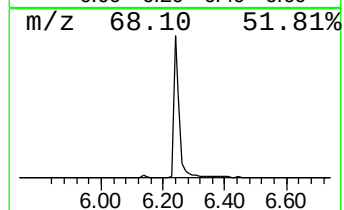
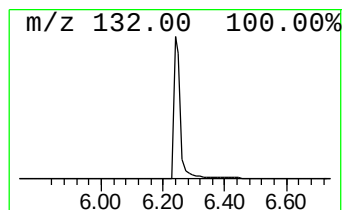
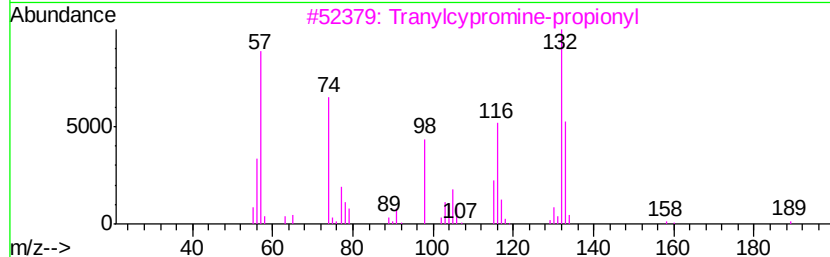
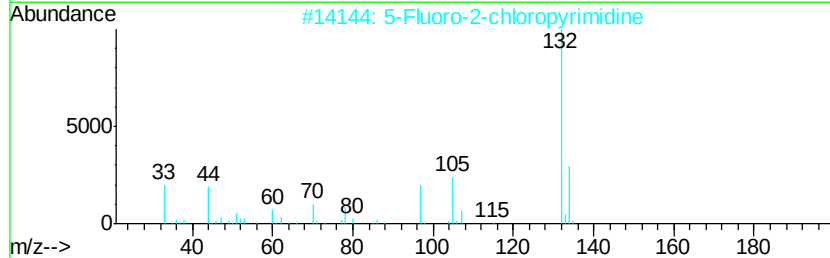
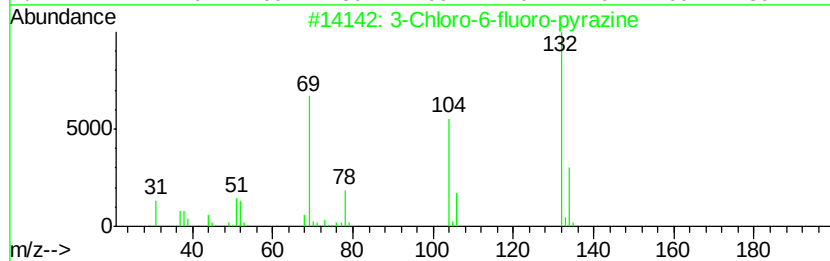
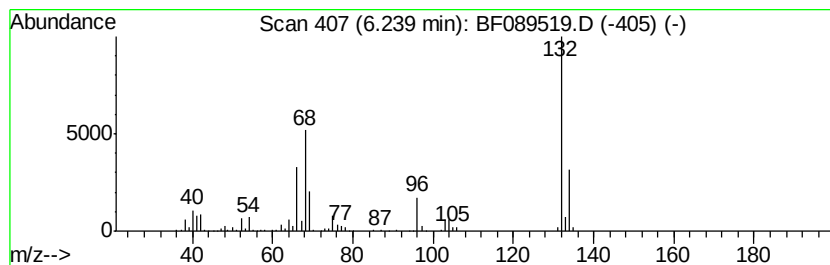
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Peak Number 3 unknown6.24 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.24	95.55 ng	970813	1,4-Dichlorobenzene-d4	6.48

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	27
2		5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
3		Tranvlcypropamine-propionyl	189	C12H15NO	1000123-86-3	10
4		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
5		Cyclopropanamine, 1-phenyl-	133	C9H11N	1000351-26-2	10



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.56	7.6	ng	76800	1	6.48	203209	20.0
unknown6.24	6.24	95.5	ng	970813	1	6.48	203209	20.0