

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF092017\
 Data File : BF098663.D
 Acq On : 20 Sep 2017 16:40
 Operator : SJ/JU
 Sample : I5276-02
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SASP2P-ENV-117-4(5-10)

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-BF091117.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	4.793	457	460	472	rBV	281434	452283	16.11%	2.026%
2	5.228	530	534	553	rBV	2652632	2732196	97.34%	12.236%
3	6.281	708	713	728	rBV	2587266	2763001	98.43%	12.374%
4	6.392	728	732	735	rBV	3111006	2806956	100.00%	12.571%
5	6.616	767	770	780	rVB	649305	594112	21.17%	2.661%
6	6.775	793	797	810	rBV	2217344	1878925	66.94%	8.415%
7	7.192	863	868	871	rBV	1804257	1657366	59.04%	7.423%
8	7.904	985	989	1006	rVB	897196	761418	27.13%	3.410%
9	8.981	1167	1172	1175	rBV	3389347	2766077	98.54%	12.388%
10	9.380	1236	1240	1247	rBV	261127	216828	7.72%	0.971%
11	9.651	1282	1286	1290	rBV	974169	787907	28.07%	3.529%
12	10.445	1417	1421	1424	rBV	1603058	1365848	48.66%	6.117%
13	11.133	1535	1538	1542	rBV	732899	644939	22.98%	2.888%
14	12.721	1803	1808	1811	rBV	1703459	1579013	56.25%	7.072%
15	13.621	1957	1961	1969	rBV3	24643	38032	1.35%	0.170%
16	13.774	1983	1987	1990	rVB	506665	503380	17.93%	2.254%
17	14.515	2110	2113	2121	rBV	61241	79052	2.82%	0.354%
18	15.168	2219	2224	2233	rBV	610323	701440	24.99%	3.141%

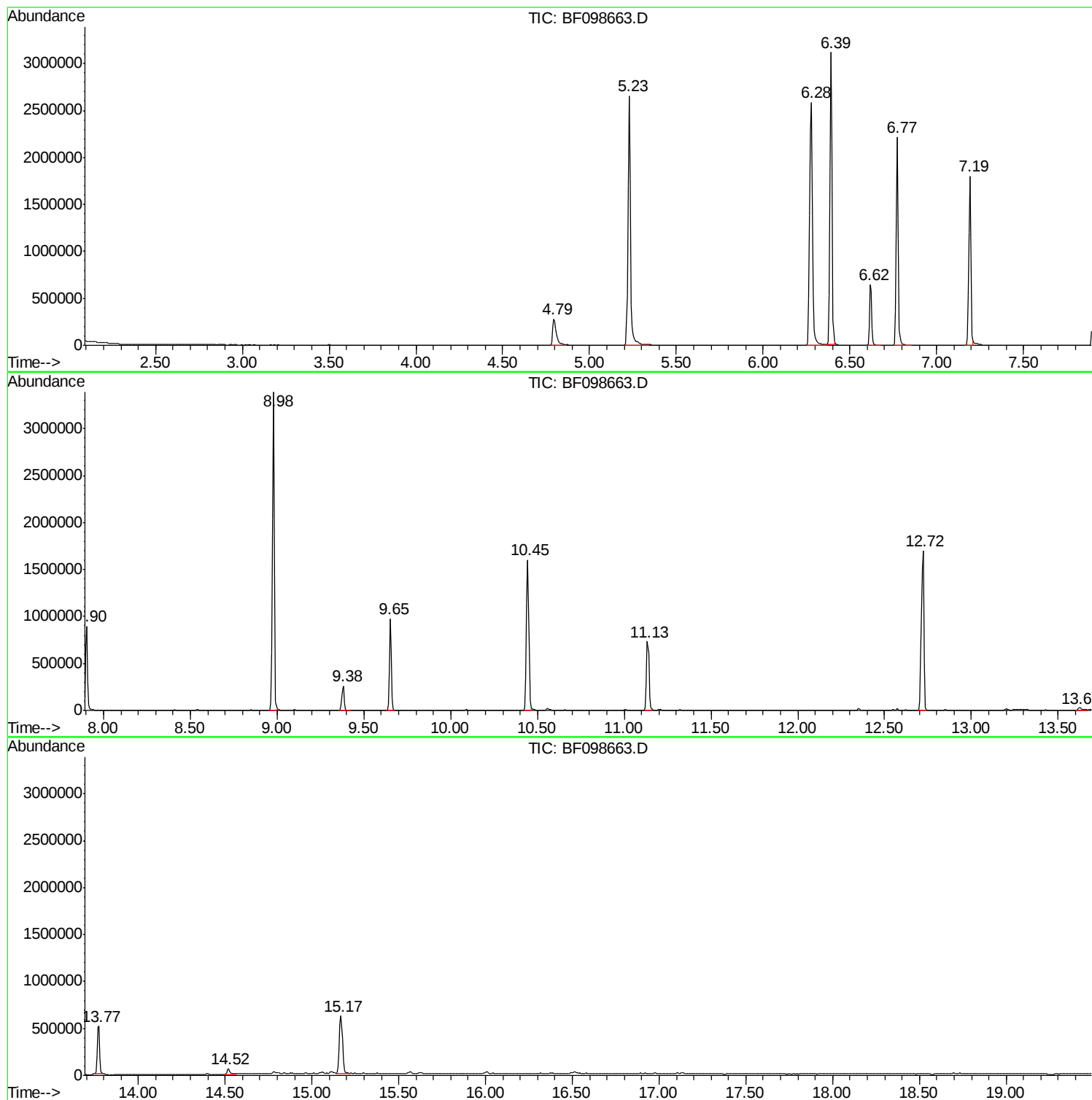
Sum of corrected areas: 22328773

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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF091117.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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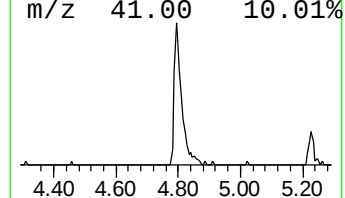
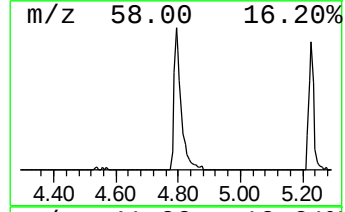
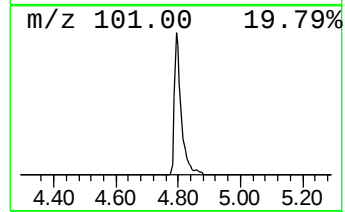
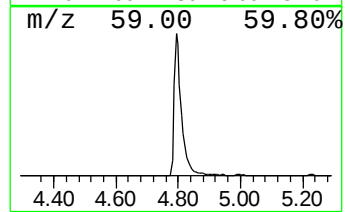
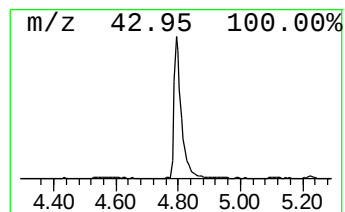
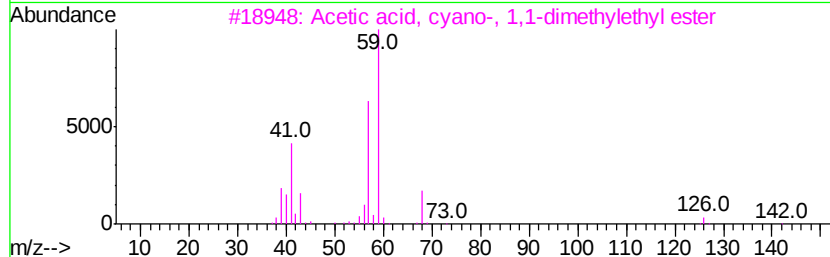
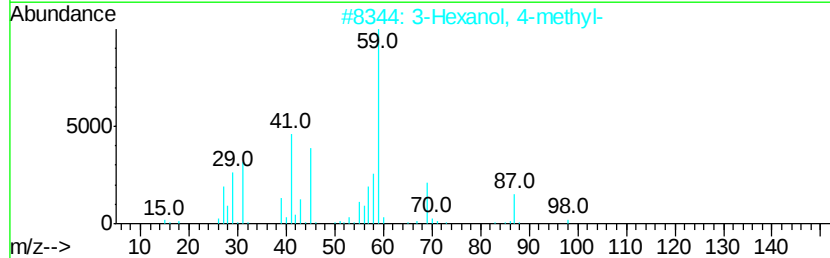
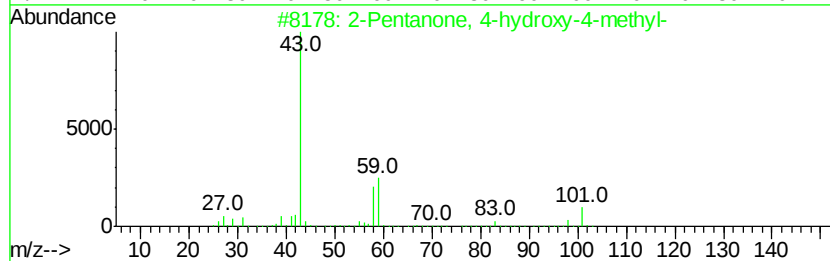
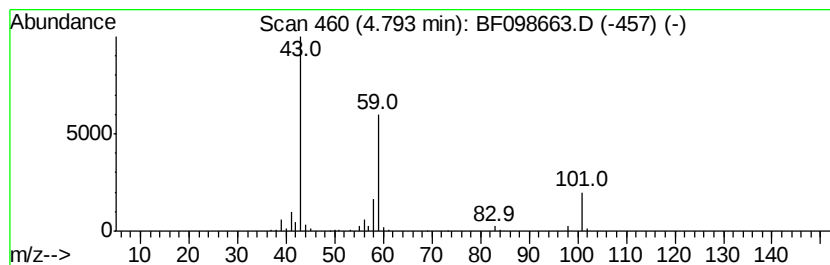
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF091117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.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.
4.79	15.23 ng	452283	1,4-Dichlorobenzene-d4	6.62

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	64		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	28		
3	Acetic acid, cyano-, 1,1-dimethyl...	141	C7H11NO2	001116-98-9	17		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9		
5	Butane, 1-ethoxy-	102	C6H14O	000628-81-9	9		



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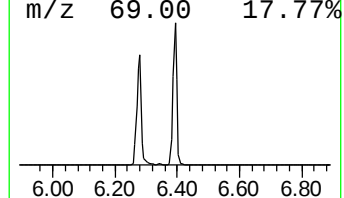
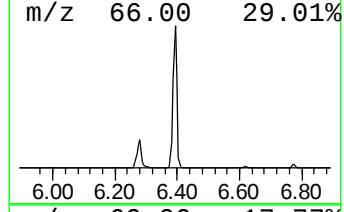
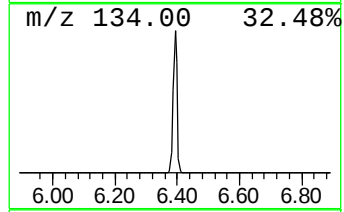
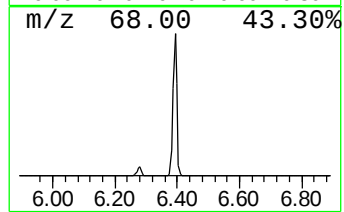
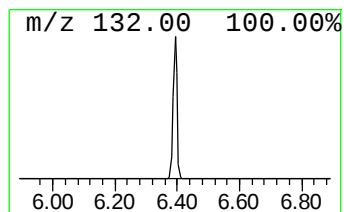
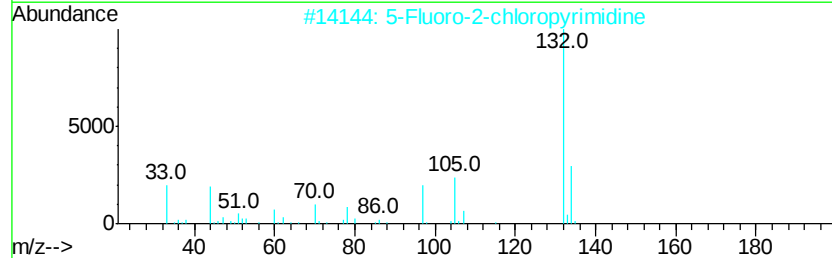
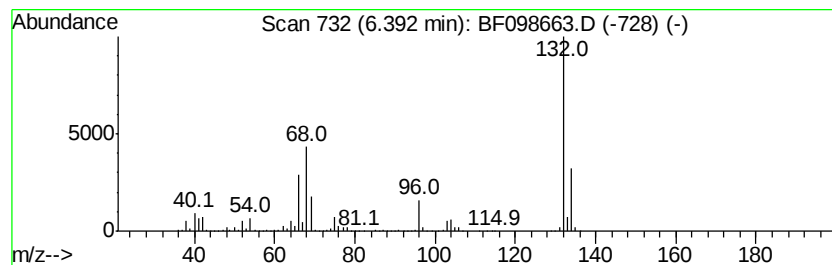
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Peak Number 2 unknown6.39 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.39	94.49 ng	2806960	1,4-Dichlorobenzene-d4	6.62

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		Tranvlcvpromine-propionyl	189	C12H15NO	1000123-86-3	10
4		Xenon	132	Xe	007440-63-3	9
5		4-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-60-2	9



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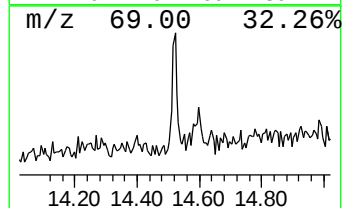
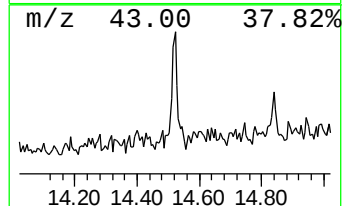
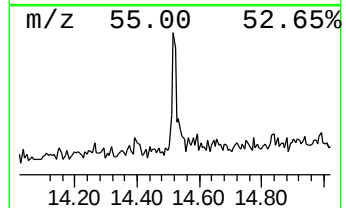
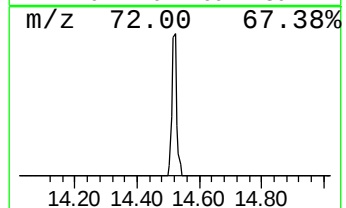
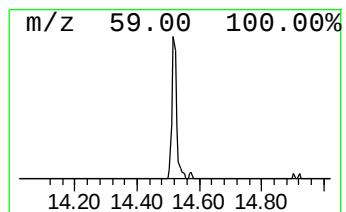
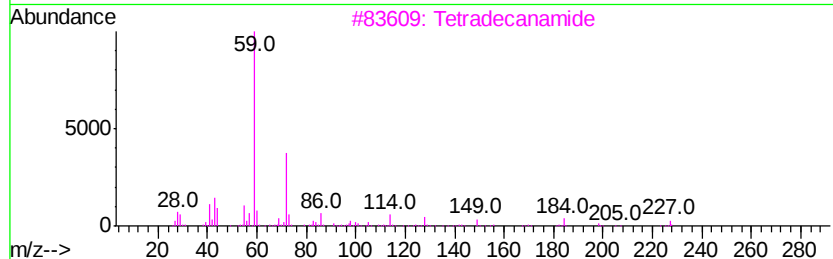
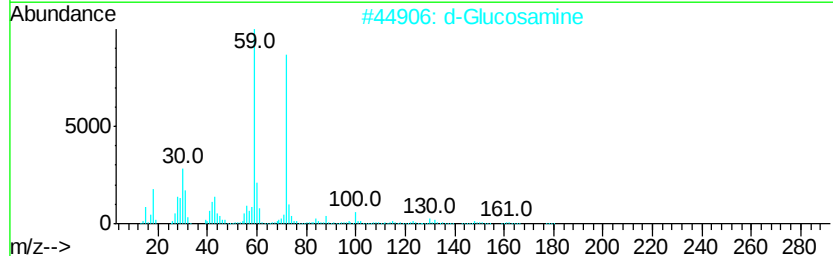
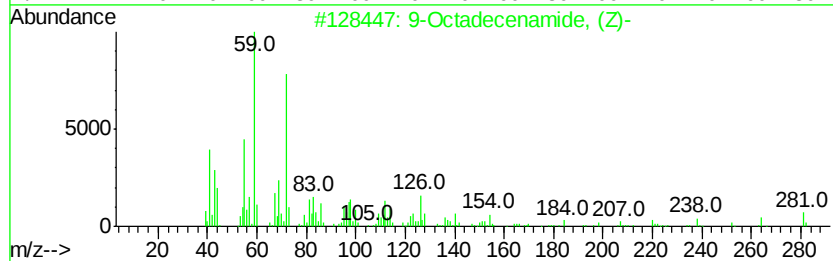
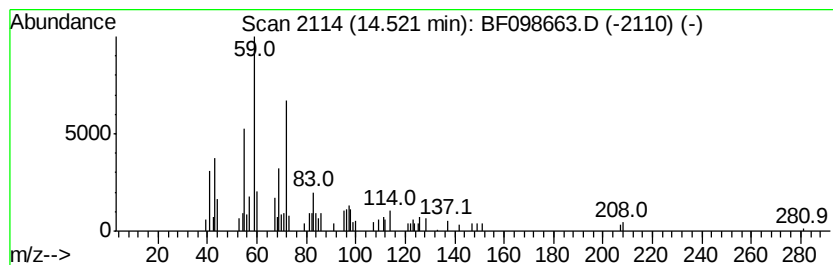
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Peak Number 4 9-Octadecenamide, (Z)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	2.25 ng	79052	Perylene-d12	15.17

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	87
2		d-Glucosamine	179	C6H13NO5	000090-77-7	53
3		Tetradecanamide	227	C14H29NO	000638-58-4	50
4		Octanamide	143	C8H17NO	000629-01-6	47
5		7-Nonenamide	155	C9H17NO	090949-53-4	47



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TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2-Pentanone, 4-hy...	4.79	15.2	ng	452283	1	6.62	594112	20.0
unknown6.39	6.39	94.5	ng	2806960	1	6.62	594112	20.0
9-Octadecenamide, ...	14.52	2.3	ng	79052	6	15.17	701440	20.0