

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF090319\
 Data File : BF116787.D
 Acq On : 3 Sep 2019 21:53
 Operator : HP/JU
 Sample : K4580-03 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 T12-B1-B4-(0-3.5)

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:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF083019.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	5.457	569	573	582	rVB	459935	417150	36.88%	4.476%
2	6.469	741	745	755	rBV	572527	471539	41.69%	5.060%
3	6.616	766	770	773	rBV	517542	478374	42.29%	5.133%
4	6.845	806	809	813	rBV	859245	709447	62.72%	7.612%
5	7.004	832	836	839	rBV	465072	369629	32.68%	3.966%
6	7.398	900	903	911	rBV	317638	273723	24.20%	2.937%
7	8.128	1024	1027	1031	rBV	1245906	972128	85.95%	10.431%
8	8.886	1153	1156	1163	rBV	18603	23387	2.07%	0.251%
9	9.198	1206	1209	1213	rBV	742462	586990	51.90%	6.298%
10	9.592	1273	1276	1282	rBV	87929	70826	6.26%	0.760%
11	9.880	1321	1325	1329	rBV	1419501	1131064	100.00%	12.136%
12	10.663	1455	1458	1462	rBV	354410	316878	28.02%	3.400%
13	10.804	1479	1482	1484	rBV	25333	21388	1.89%	0.229%
14	11.363	1574	1577	1580	rBV	1217989	1063722	94.05%	11.414%
15	12.945	1843	1846	1850	rVB	539933	482568	42.66%	5.178%
16	13.998	2022	2025	2029	rVB	890813	796680	70.44%	8.548%
17	14.515	2111	2113	2119	rVB	269927	304880	26.96%	3.271%
18	15.457	2268	2273	2276	rBV	679941	829465	73.33%	8.900%

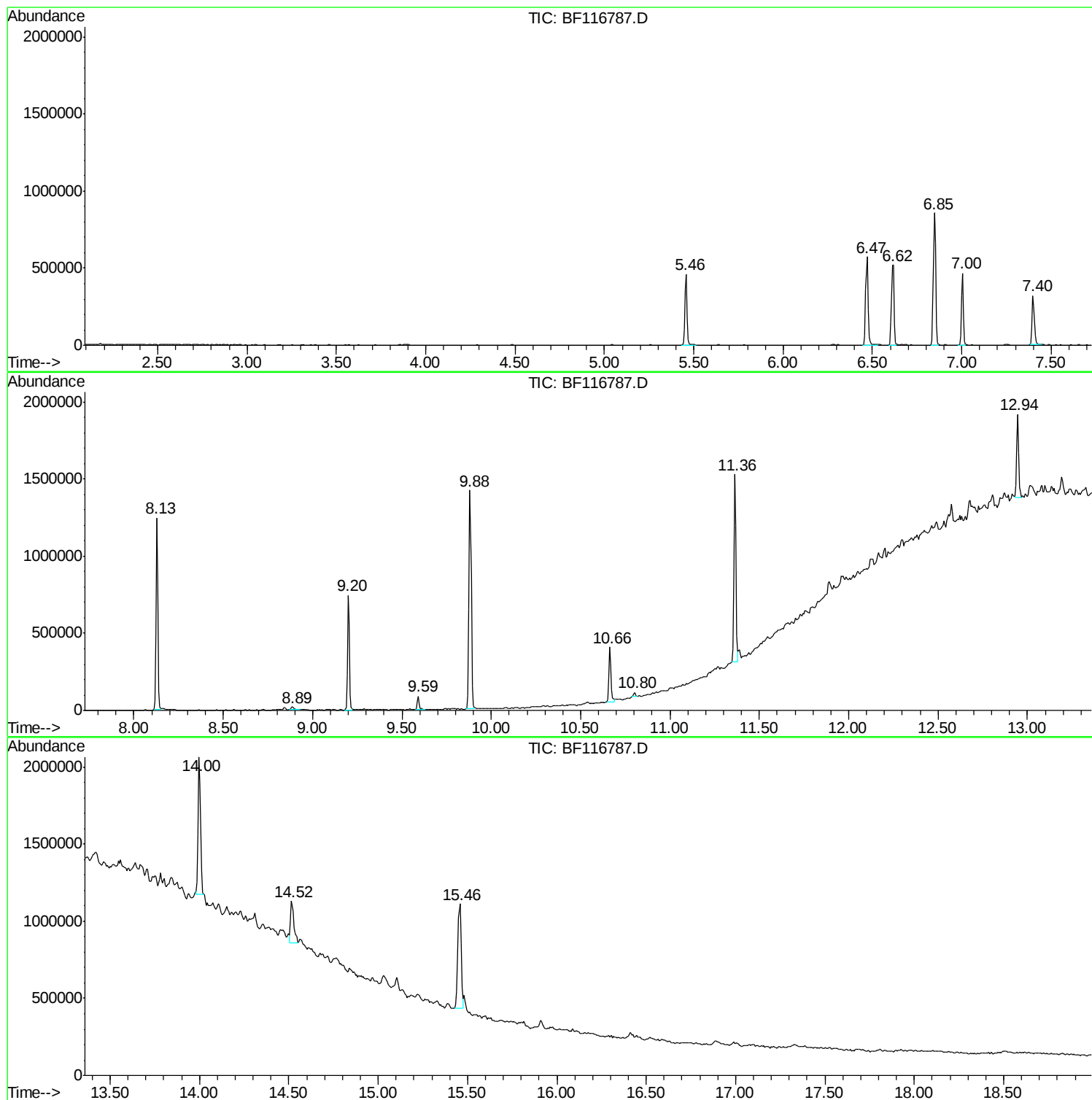
Sum of corrected areas: 9319838

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.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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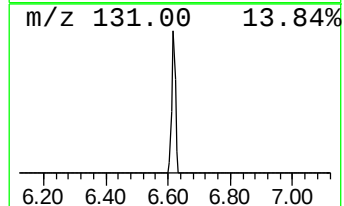
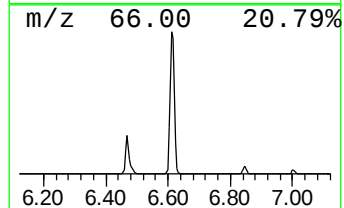
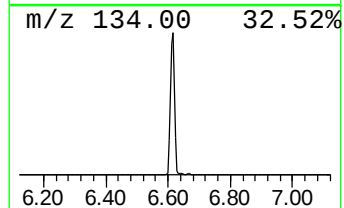
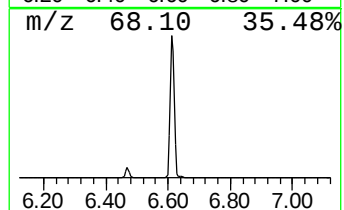
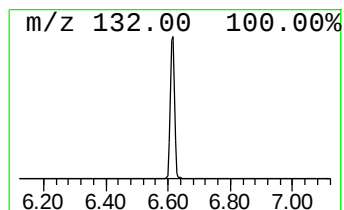
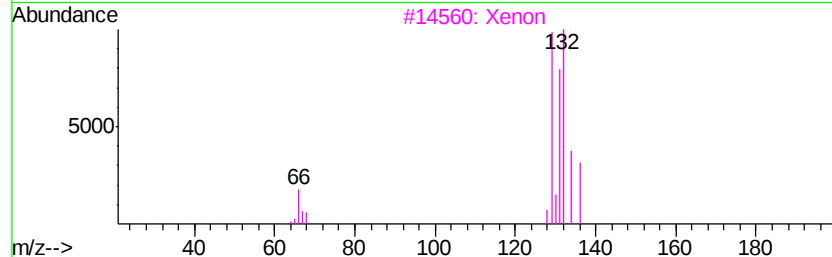
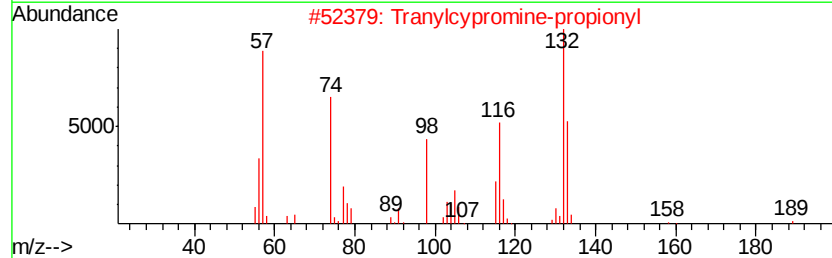
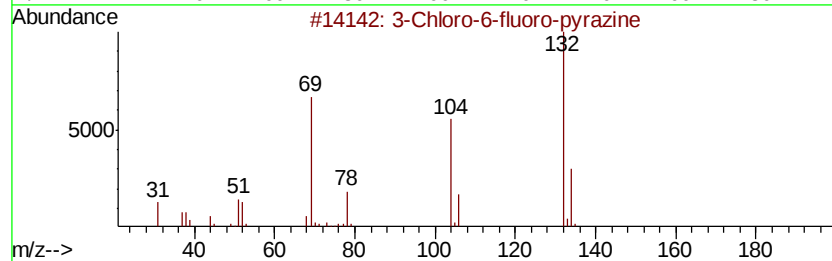
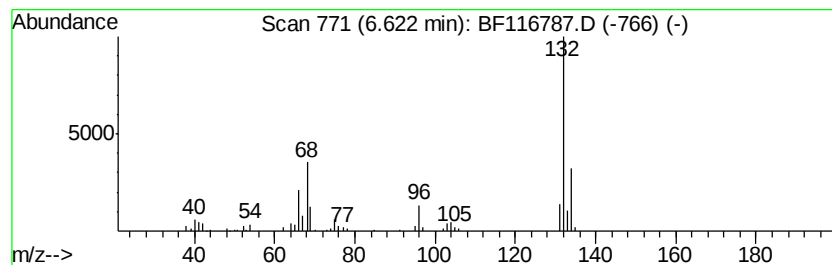
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
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TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown6.62 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.62	13.49 ng	478374	1,4-Dichlorobenzene-d4	6.85

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Chloro-6-fluoro-pyrazine	132	C4H2ClFN2	1000146-10-7	25
2		Tranvlcypropromine-propionyl	189	C12H15NO	1000123-86-3	16
3		Xenon	132	Xe	007440-63-3	9
4		1-Methylpyrrolo[1,2-a]pyrazine	132	C8H8N2	064608-59-9	9
5		2H-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	9



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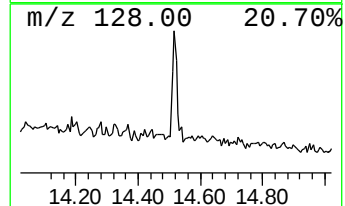
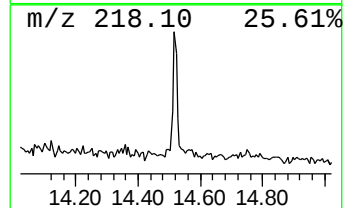
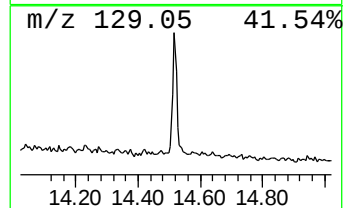
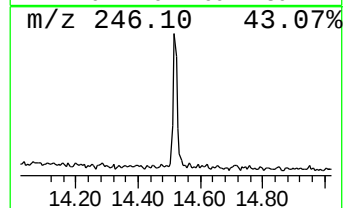
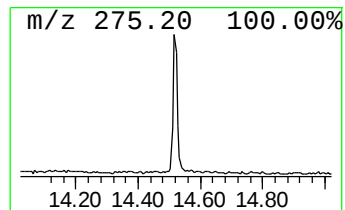
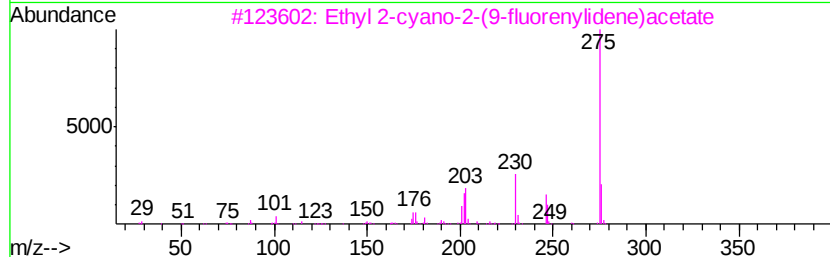
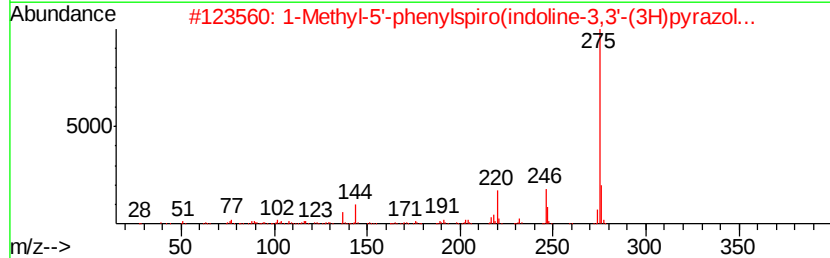
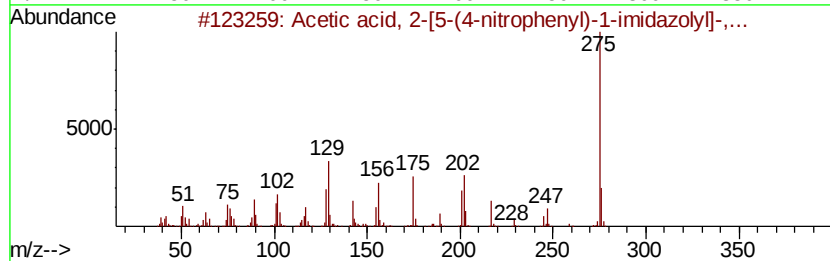
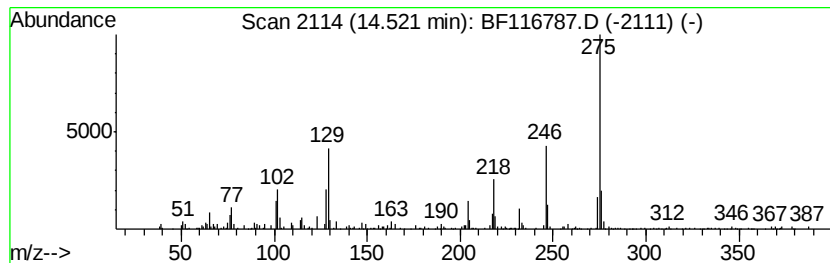
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Peak Number 3 unknown14.52 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.52	7.65 ng	304880	Chrysene-d12	14.00

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Acetic acid, 2-[5-(4-nitrophenyl)-1-imidazolyl]-,...	275	C13H13N3O4	351064-45-4	49
2		1-Methyl-5'-phenylspiro(indoline-3,3'-(3H)pyrazol...	275	C17H13N3O	015970-21-5	38
3		Ethyl 2-cyano-2-(9-fluorenylidene)acetate	275	C18H13N02	014003-25-9	35
4		8H-Benzo[ql]-1,3-benzodioxolo[6,5-b]pyridine	275	C17H9N03	000475-75-2	30
5		1,4-Benzenediamine, N,N-dimethyl-2,5-bis(4-chlorophenyl)-	275	C13H13N3O2S	126893-36-5	27



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.62	6.62	13.5	ng	478374	1	6.85	709447	20.0
unknown14.52	14.52	7.7	ng	304880	5	14.00	796680	20.0