

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF061219\
 Data File : BF114975.D
 Acq On : 12 Jun 2019 15:31
 Operator : HP/JU
 Sample : PB120267BL
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB120267BL

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-BF061219.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	5.269	535	541	544	rBV	5642506	6201690	76.25%	9.718%
2	6.298	709	716	718	rBV	5271945	6319047	77.69%	9.902%
3	6.422	731	737	740	rBV	6059332	6632754	81.55%	10.393%
4	6.645	771	775	779	rBV	1958945	1579219	19.42%	2.475%
5	6.804	797	802	805	rBV	5684528	5408340	66.49%	8.475%
6	7.210	865	871	874	rBV	3759235	4298051	52.84%	6.735%
7	7.928	988	993	996	rBV	2181926	2057565	25.30%	3.224%
8	9.010	1170	1177	1180	rBV	6657573	8133458	100.00%	12.745%
9	9.675	1284	1290	1294	rBV	2658732	2415561	29.70%	3.785%
10	10.463	1417	1424	1427	rBV	4803247	5017005	61.68%	7.861%
11	11.151	1535	1541	1544	rBV	2859850	2591479	31.86%	4.061%
12	12.739	1805	1811	1814	rBV	7234986	8095264	99.53%	12.685%
13	13.768	1982	1986	1990	rBV	2388088	2343020	28.81%	3.671%
14	15.156	2217	2222	2228	rBV	1792953	2047726	25.18%	3.209%
15	15.804	2321	2332	2348	rVB3	169567	677506	8.33%	1.062%

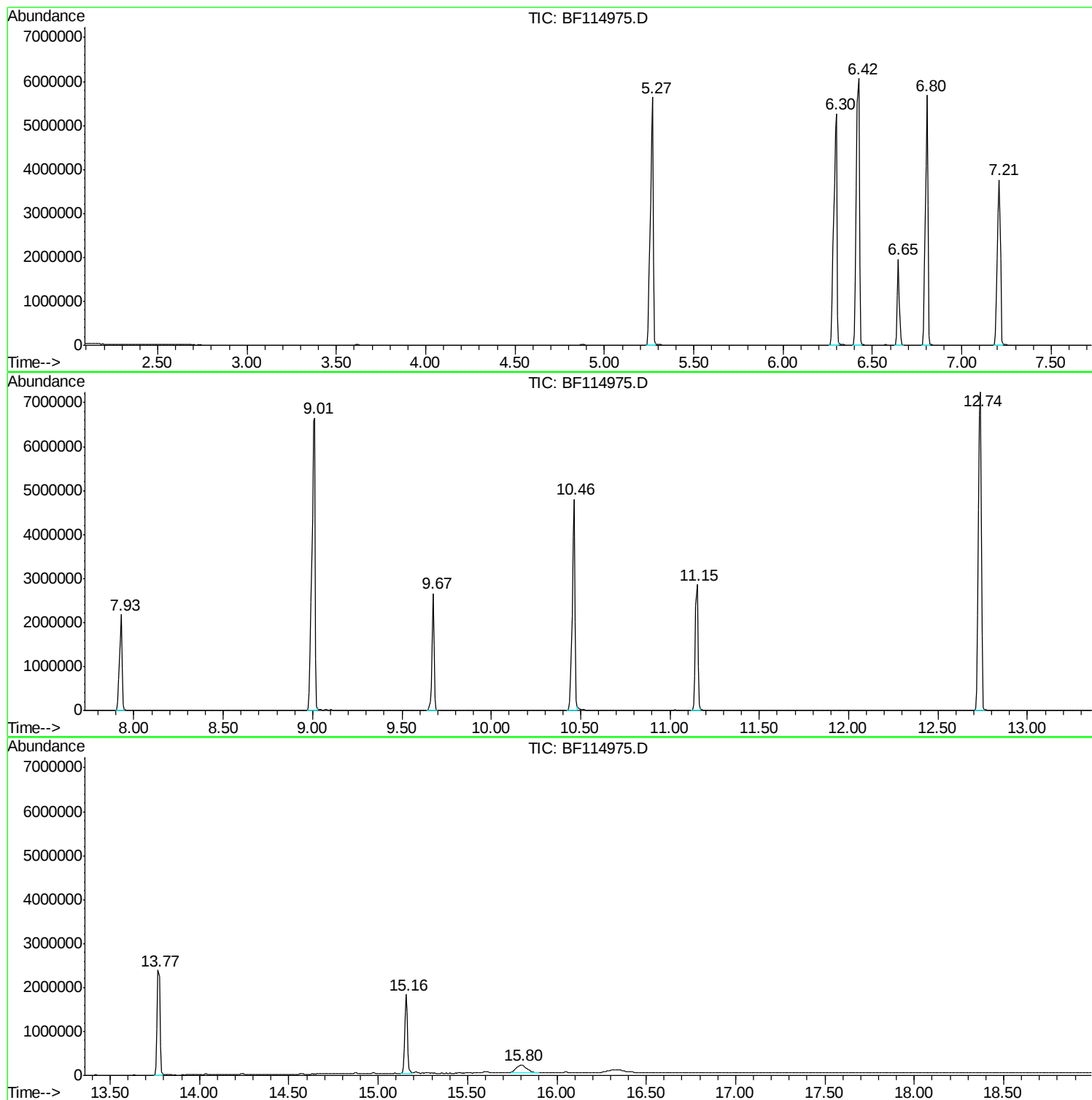
Sum of corrected areas: 63817685

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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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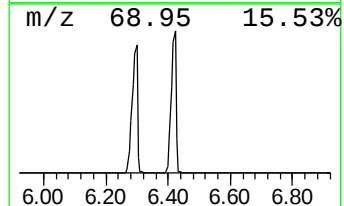
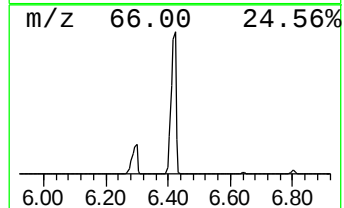
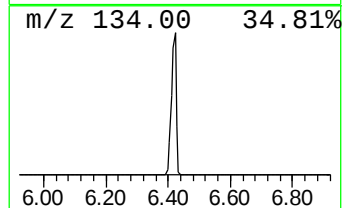
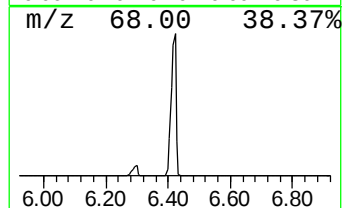
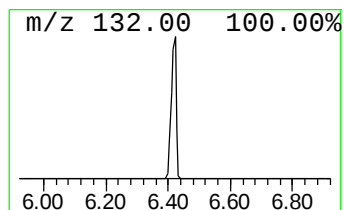
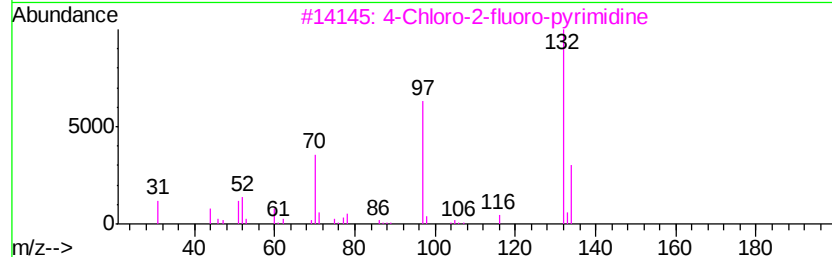
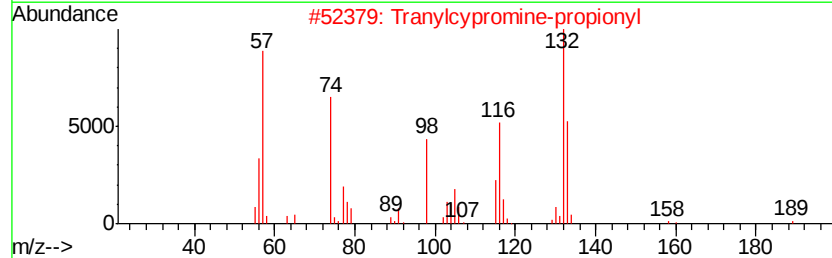
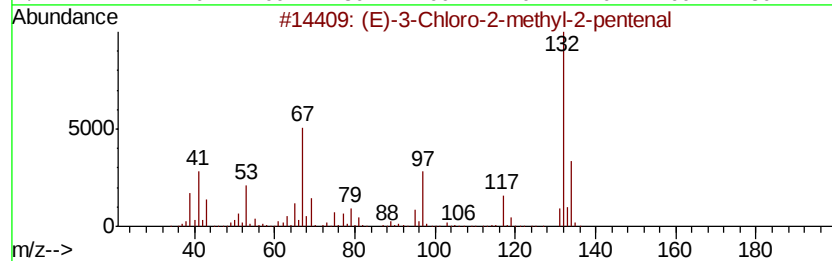
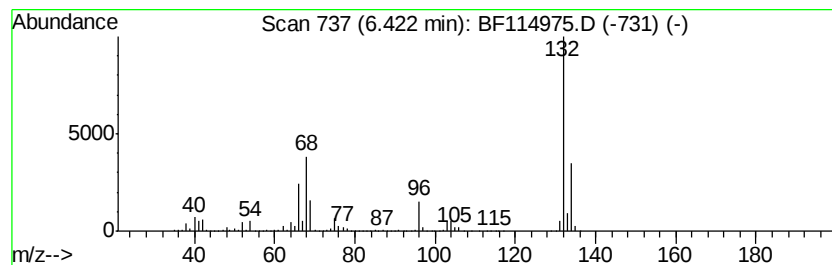
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Peak Number 1 unknown6.42 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.42	84.00 ng	6632750	1,4-Dichlorobenzene-d4	6.65

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	16
2		Tranylcypromine-propionyl	189	C12H15NO	1000123-86-3	12
3		4-Chloro-2-fluoro-pyrimidine	132	C4H2ClFN2	051422-00-5	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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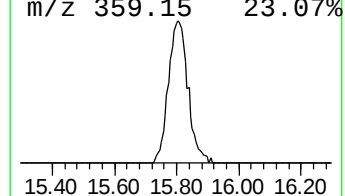
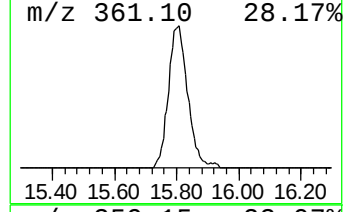
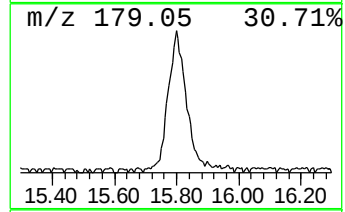
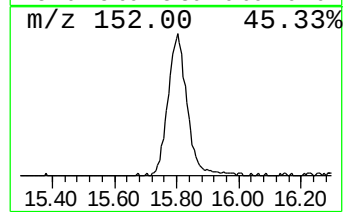
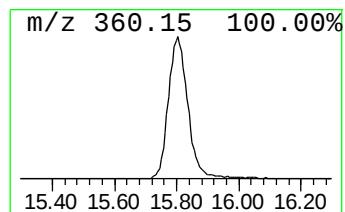
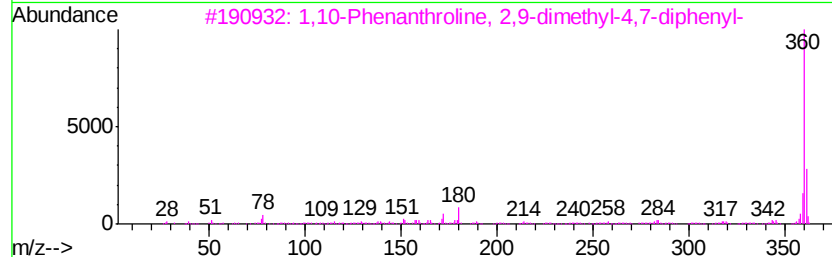
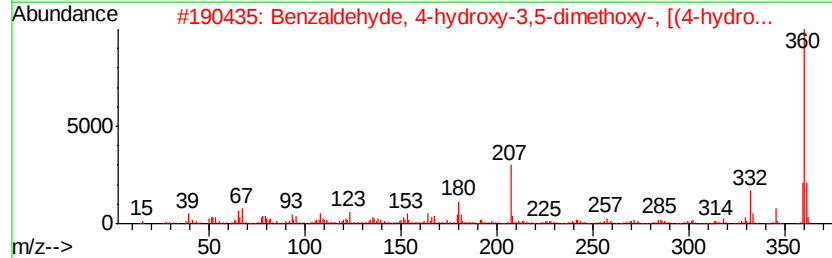
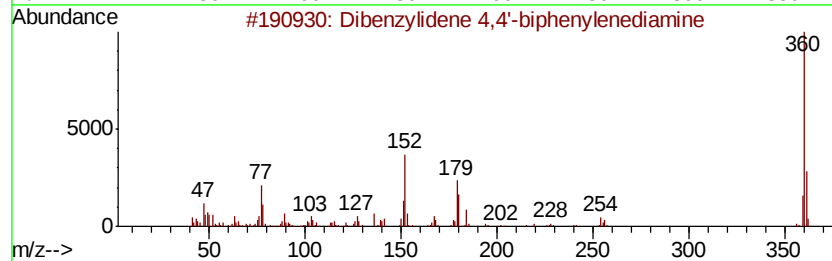
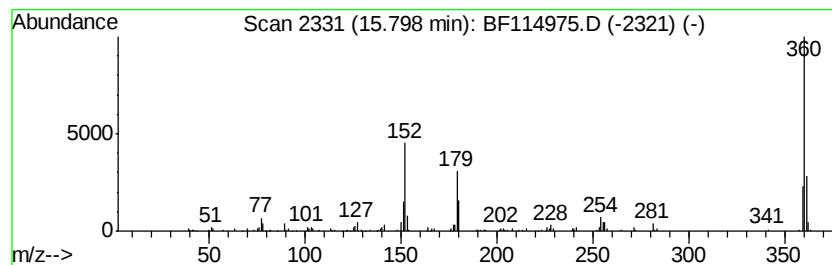
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Peak Number 3 Dibenzylidene 4,4'-biphenyl... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.80	6.62 ng	677506	Perylene-d12	15.16

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Dibenzylidene 4,4'-biphenylenedi...	360	C26H20N2	006311-48-4	94
2		Benzaldehyde, 4-hydroxy-3,5-dime...	360	C18H20N2O6	014414-32-5	49
3		1,10-Phenanthroline, 2,9-dimethv...	360	C26H20N2	004733-39-5	49
4		N,N'-di-2-Naphthyl-p-phenylenedi...	360	C26H20N2	000093-46-9	46
5		1,3,2-Dioxaborolane, bis(spiro[4...	360	C20H13B06	1000150-23-7	43



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown6.42	6.42	84.0	ng	6632750	1	6.65	1579220	20.0
Dibenzylidene 4,4...	15.80	6.6	ng	677506	6	15.16	2047730	20.0