

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF031920\
 Data File : BF119447.D
 Acq On : 19 Mar 2020 19:27
 Operator : CG/JU
 Sample : PB127678BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB127678BL

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-BF031820.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	5.075	502	508	514	rBV	83059	112847	1.29%	0.181%
2	5.463	569	574	577	rBV	7678253	7784769	89.17%	12.483%
3	6.469	739	745	748	rBV	7178735	7537008	86.33%	12.085%
4	6.845	806	809	813	rVB	2181016	1800478	20.62%	2.887%
5	7.004	832	836	839	rBV	6803469	6102905	69.90%	9.786%
6	7.410	899	905	908	rBV	5034958	4980495	57.05%	7.986%
7	8.133	1023	1028	1031	rBV	2598403	2402202	27.52%	3.852%
8	9.216	1206	1212	1215	rBV	8110580	8730452	100.00%	13.999%
9	9.892	1321	1327	1330	rBV	3038822	2666603	30.54%	4.276%
10	10.680	1454	1461	1464	rBV	5585134	5172930	59.25%	8.295%
11	11.380	1574	1580	1583	rBV	3140045	2683925	30.74%	4.304%
12	12.974	1845	1851	1854	rBV	8080506	7858777	90.02%	12.601%
13	13.833	1991	1997	2001	rBV	195334	245393	2.81%	0.393%
14	14.021	2025	2029	2032	rBV	2032430	1845163	21.13%	2.959%
15	14.651	2131	2136	2139	rBV	332753	455069	5.21%	0.730%
16	14.674	2139	2140	2144	rVV	165865	129032	1.48%	0.207%
17	14.721	2145	2148	2155	rVB2	78967	96755	1.11%	0.155%
18	15.486	2273	2278	2284	rBV	1304285	1672360	19.16%	2.682%
19	15.986	2359	2363	2371	rVB3	53017	87507	1.00%	0.140%

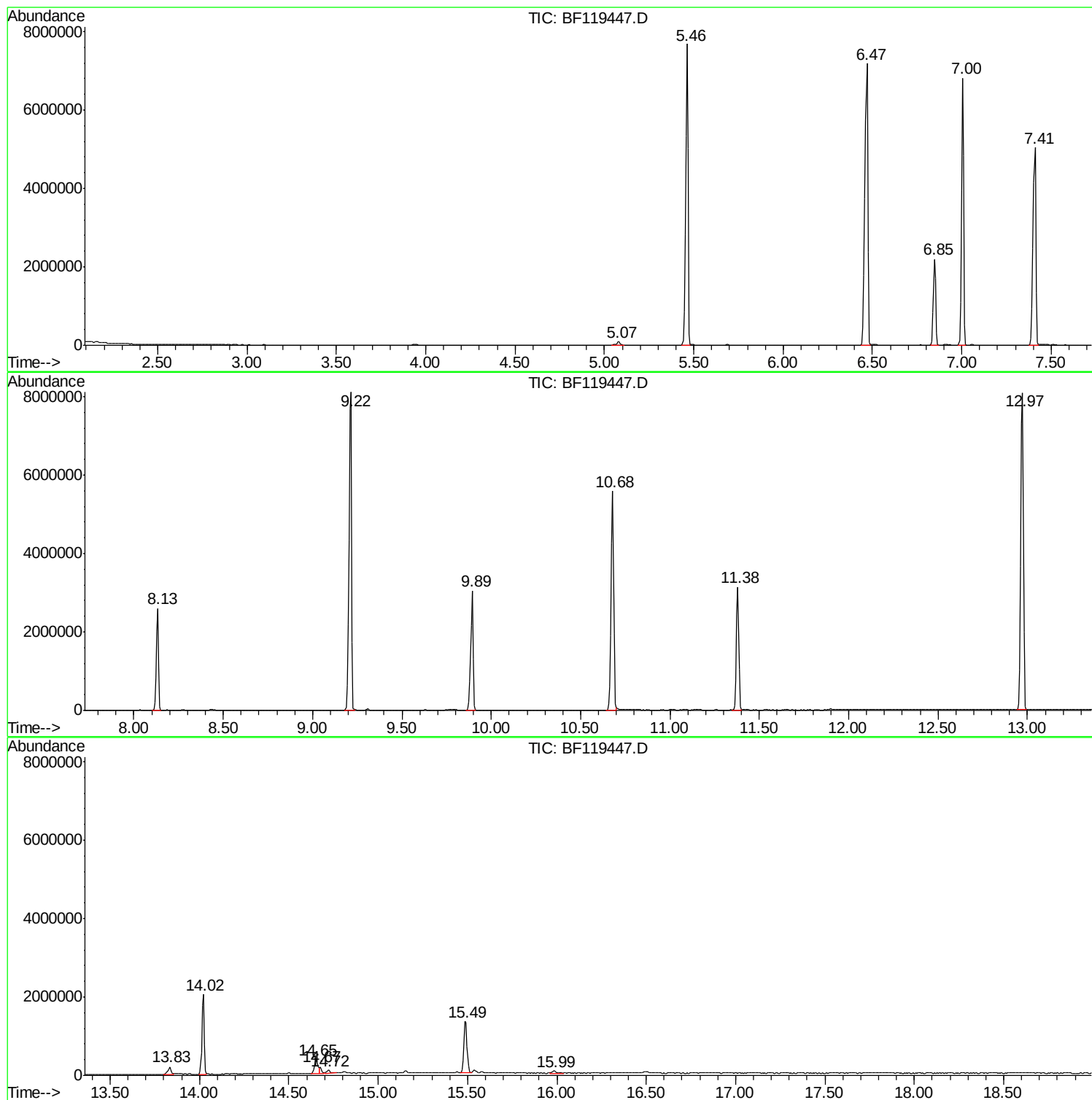
Sum of corrected areas: 62364670

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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF031820.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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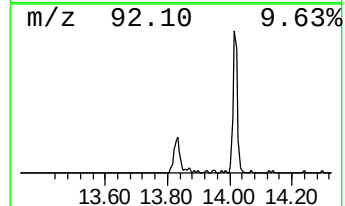
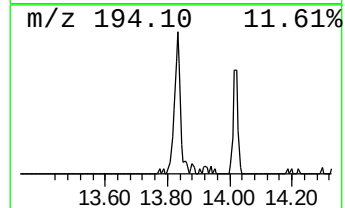
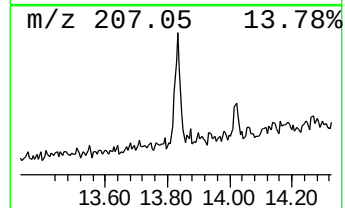
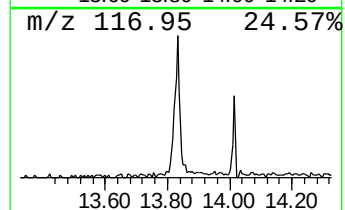
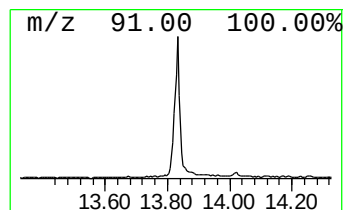
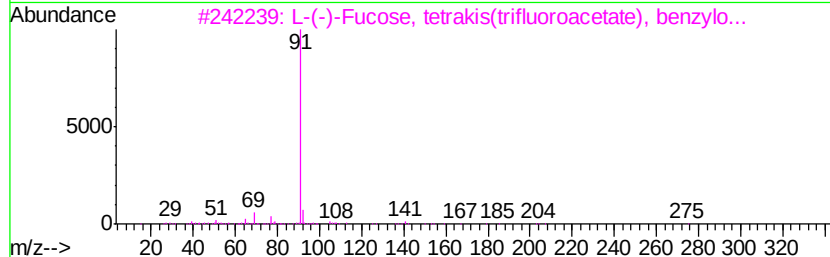
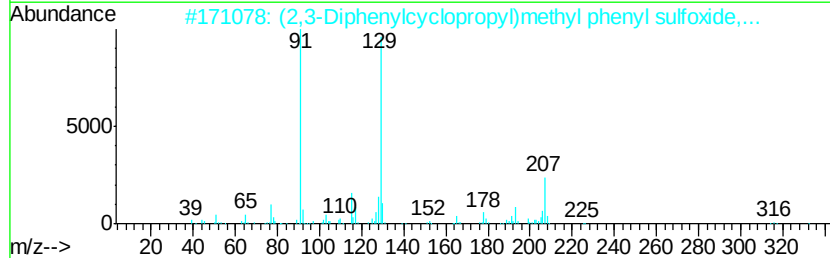
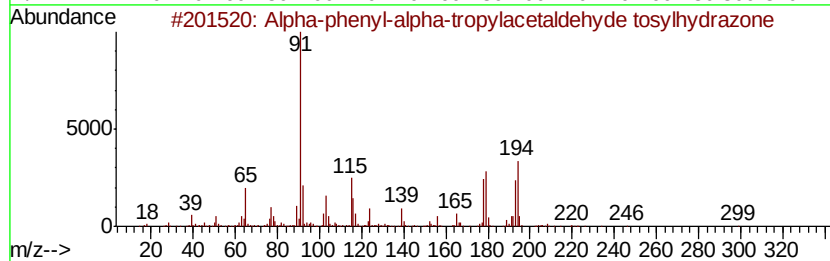
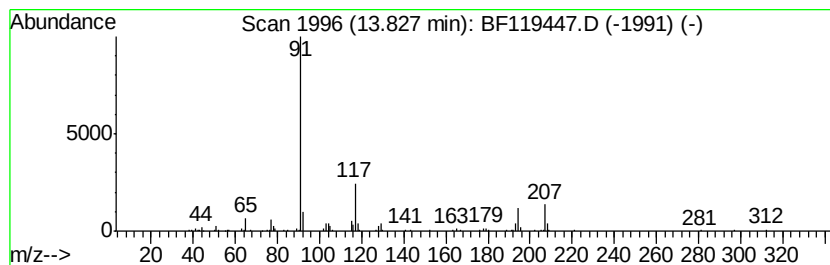
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Peak Number 2 unknown13.83 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	2.66 ng	245393	Chrysene-d12	14.02

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Alpha-phenyl-alpha-tropylacetalde...	378	C22H22N2O2S	022532-16-7	25
2		(2,3-Diphenylcyclopropyl)methyl ...	332	C22H20OS	131758-71-9	25
3		L-(-)-Fucose, tetrakis(trifluoro...	653	C21H15F12NO9	1000380-45-2	14
4		Benzene, (5-iodopentyl)-	274	C11H15I	099858-37-4	12
5		1-Benzyl-2-(trifluoromethyl)azir...	201	C10H10F3N	106240-89-5	10



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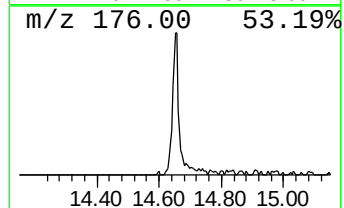
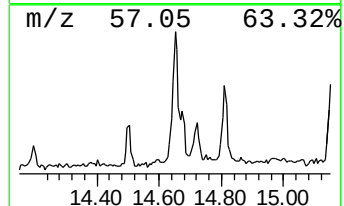
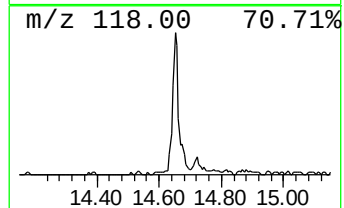
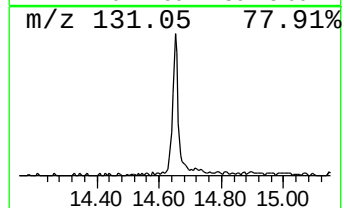
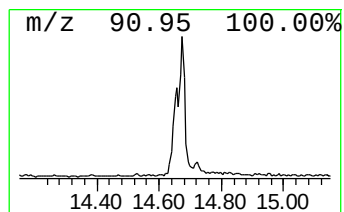
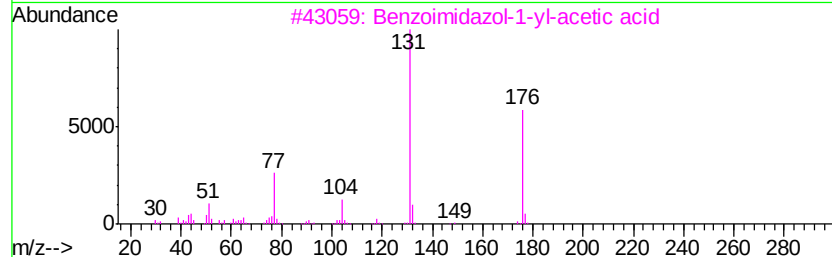
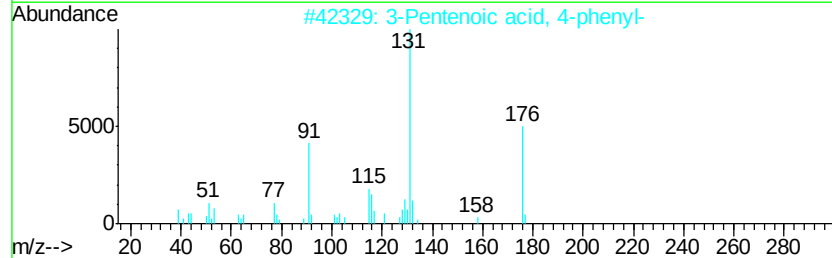
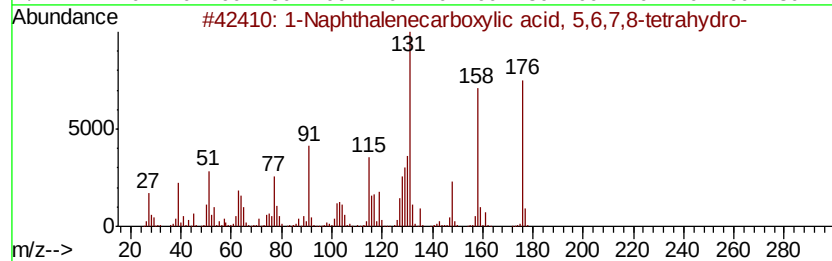
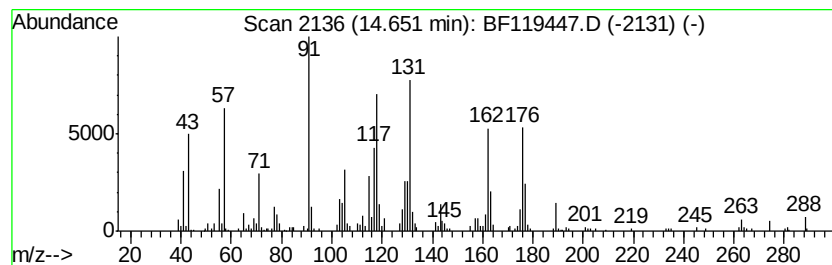
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Peak Number 3 unknown14.65 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.65	4.93 ng	455069	Chrysene-d12	14.02	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenecarboxylic acid, 5,...	176	C11H12O2	004242-18-6	43
2	3-Pentenoic acid, 4-phenyl-	176	C11H12O2	053774-19-9	38
3	Benzoimidazol-1-yl-acetic acid	176	C9H8N2O2	1000317-79-2	27
4	1-Penten-3-one, 4,4-dimethyl-1-p...	188	C13H16O	000538-44-3	25
5	.alpha.-Methyl-.alpha.-phenylsuc...	189	C11H11NO2	001497-17-2	25



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown13.83	13.83	2.7	ng	245393	5	14.02	1845160	20.0
unknown14.65	14.65	4.9	ng	455069	5	14.02	1845160	20.0