

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF122618\
 Data File : BF111740.D
 Acq On : 27 Dec 2018 3:23
 Operator : JU/SJ
 Sample : J6188-05 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-122118

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF121918.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	2.193	14	18	24	rVB	68075	90160	5.98%	0.566%
2	5.116	511	515	520	rVB	75832	74423	4.94%	0.467%
3	5.534	582	586	590	rBV	1492796	1387332	92.03%	8.713%
4	6.539	753	757	764	rBV	1631154	1507477	100.00%	9.468%
5	6.675	776	780	784	rBV	1553982	1440343	95.55%	9.046%
6	6.904	816	819	823	rVB	1013516	796636	52.85%	5.003%
7	7.063	842	846	849	rBV	1346472	1094419	72.60%	6.873%
8	7.457	909	913	917	rBV	961361	800692	53.11%	5.029%
9	8.186	1033	1037	1040	rBV	1117234	902220	59.85%	5.666%
10	8.210	1040	1041	1045	rVB	35179	20299	1.35%	0.127%
11	9.257	1216	1219	1222	rBV	1649077	1268414	84.14%	7.966%
12	9.645	1283	1285	1291	rBV	139554	129490	8.59%	0.813%
13	9.939	1332	1335	1339	rBV	1058214	932840	61.88%	5.859%
14	9.974	1339	1341	1344	rVB	41631	33294	2.21%	0.209%
15	10.145	1366	1370	1373	rVB	18389	18203	1.21%	0.114%
16	10.486	1425	1428	1431	rVB	34364	29974	1.99%	0.188%
17	10.727	1465	1469	1473	rBV	1095846	894597	59.34%	5.618%
18	10.845	1486	1489	1494	rVB5	12812	23442	1.56%	0.147%
19	11.321	1568	1570	1574	rVB3	16274	18114	1.20%	0.114%
20	11.427	1584	1588	1590	rBV	1216535	994697	65.98%	6.247%
21	11.451	1590	1592	1595	rVV	225477	169537	11.25%	1.065%
22	11.504	1598	1601	1604	rVB	59243	55358	3.67%	0.348%
23	11.651	1622	1626	1629	rBV2	30433	35469	2.35%	0.223%
24	11.933	1671	1674	1675	rBV	19001	18383	1.22%	0.115%
25	11.957	1675	1678	1682	rVB2	44381	53348	3.54%	0.335%
26	12.039	1689	1692	1695	rBV	40119	43639	2.89%	0.274%
27	12.498	1768	1770	1772	rBV3	25605	25115	1.67%	0.158%
28	12.521	1772	1774	1778	rVB2	56760	49281	3.27%	0.310%
29	12.639	1791	1794	1798	rBV	273326	217356	14.42%	1.365%
30	12.686	1799	1802	1803	rBV3	27873	24879	1.65%	0.156%
31	12.868	1828	1833	1838	rBV	221060	297853	19.76%	1.871%
32	13.010	1853	1857	1860	rVB	1436246	1194408	79.23%	7.501%
33	14.062	2033	2036	2039	rBV	793130	795258	52.75%	4.995%
34	15.551	2286	2289	2293	rBV	415939	485529	32.21%	3.049%

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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-BF121918.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

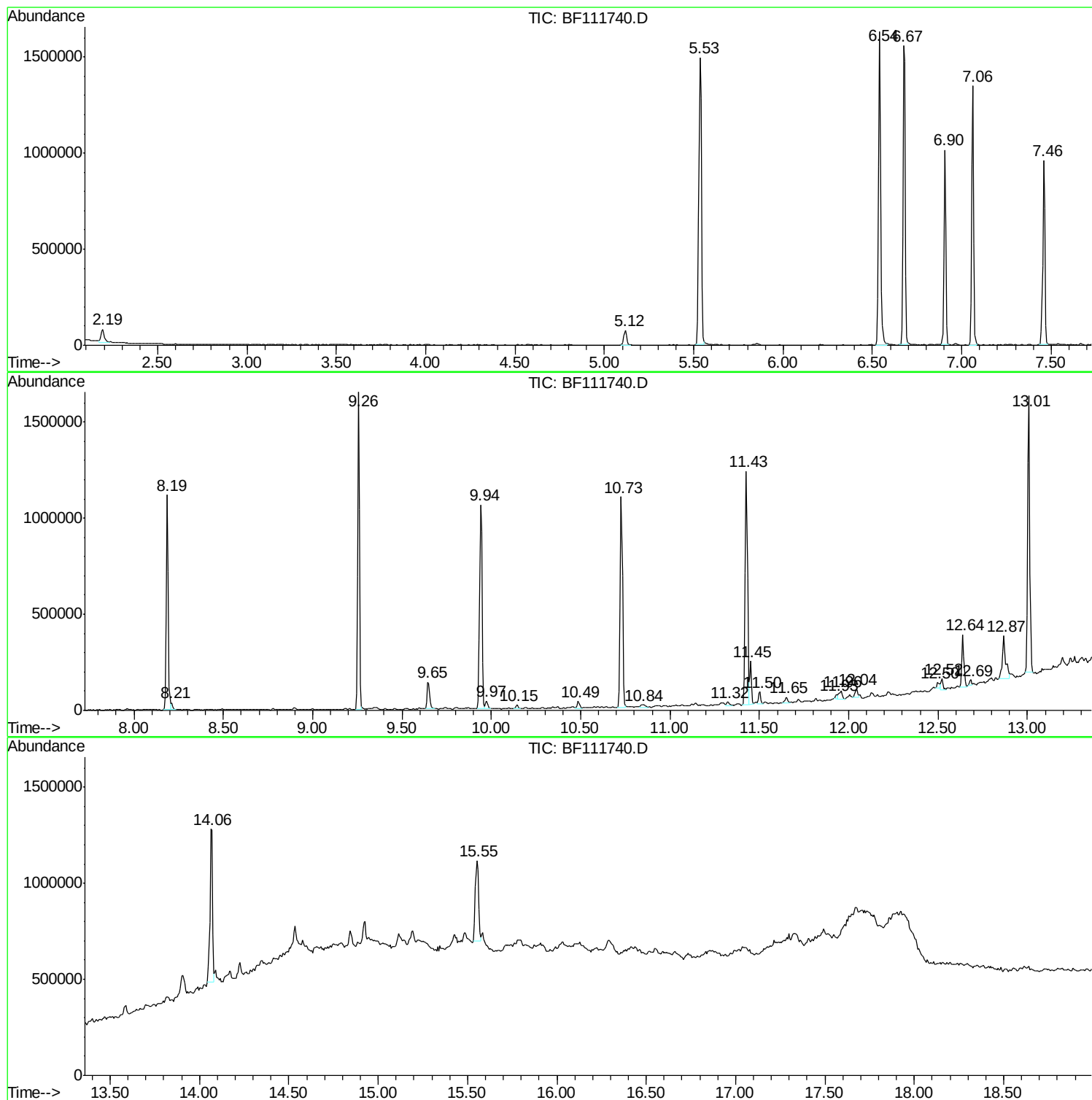
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ClientSampled :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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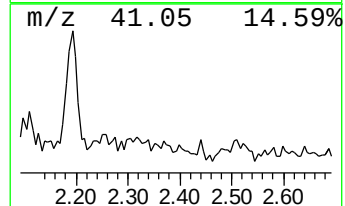
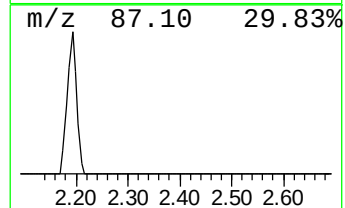
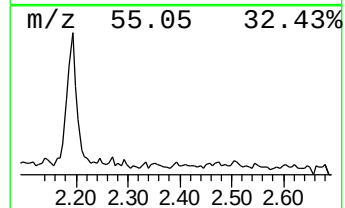
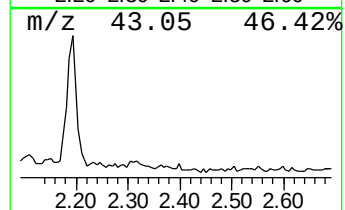
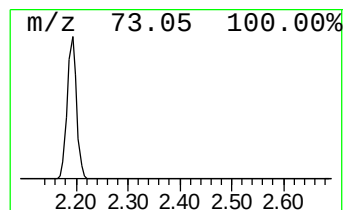
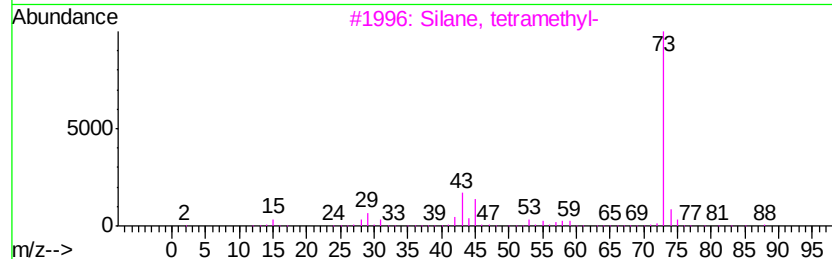
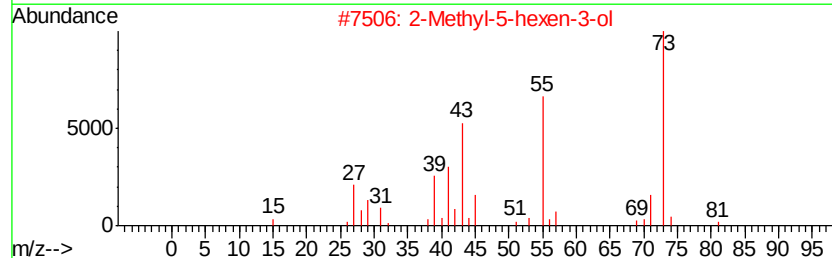
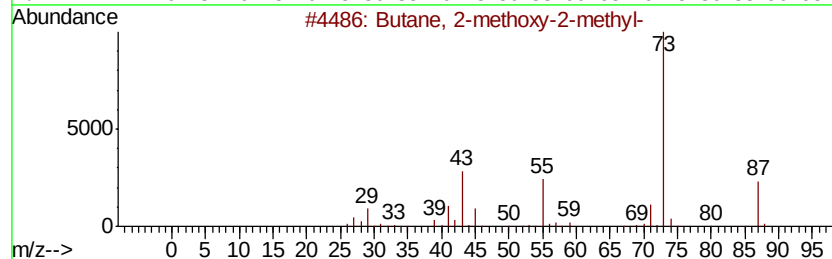
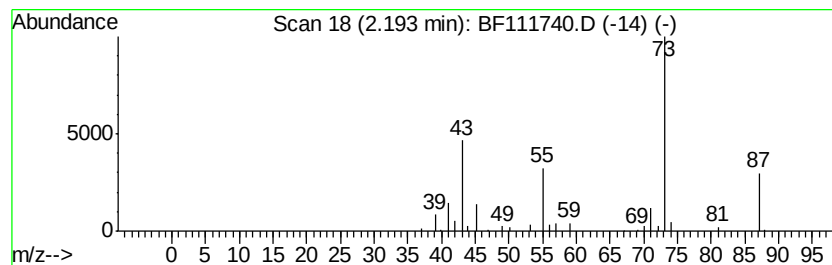
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Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.19	2.26 ng	90160	1,4-Dichlorobenzene-d4	6.90

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2-Methyl-5-hexen-3-ol	114	C7H14O	032815-70-6	37
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	33
5			Boronic acid, ethyl-, dimethyl e...	102	C4H11BO2	007318-82-3	32



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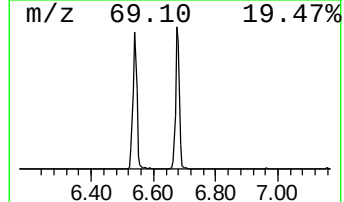
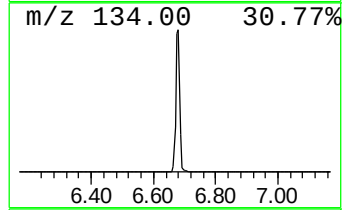
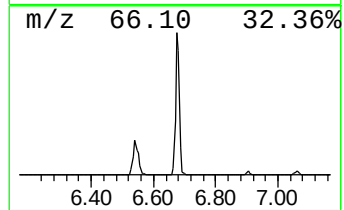
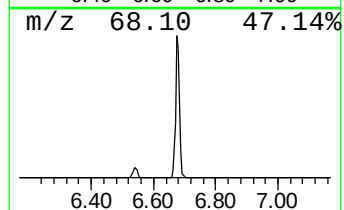
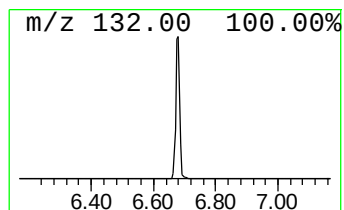
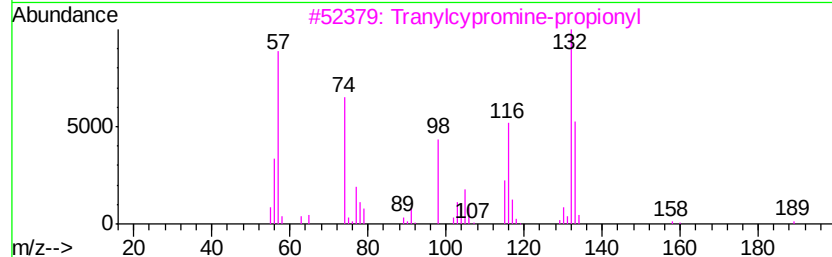
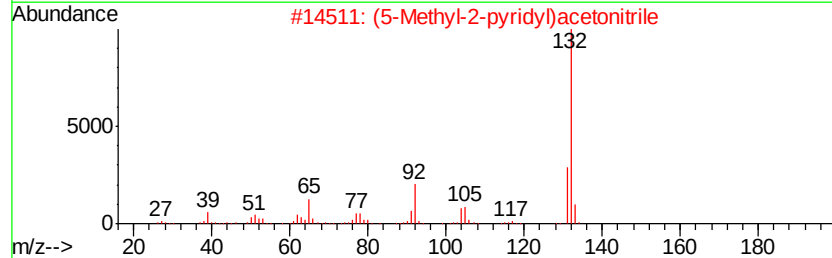
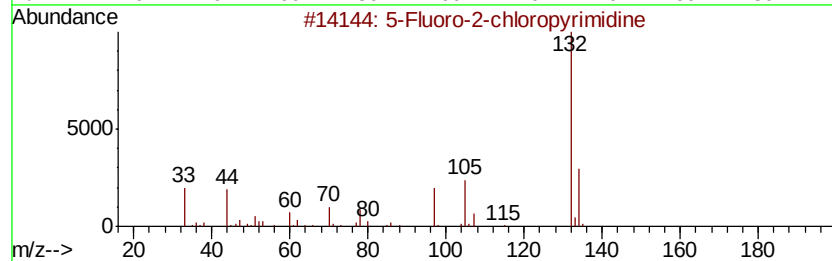
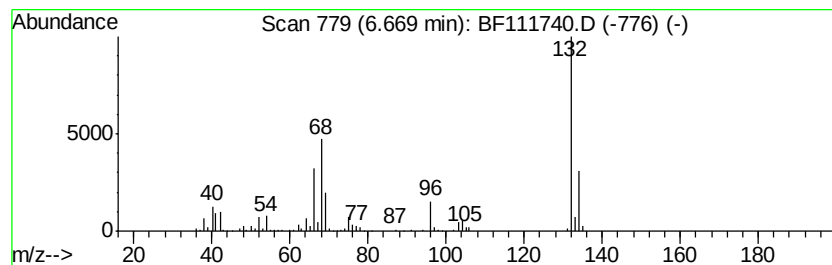
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Peak Number 2 unknown6.67 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.67	36.16 ng	1440340	1,4-Dichlorobenzene-d4	6.90

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	5-Fluoro-2-chloropyrimidine	132	C4H2ClFN2	062802-42-0	12
2		(5-Methyl-2-pyridyl)acetonitrile	132	C8H8N2	1000241-93-9	10
3		Tranvlcyproline-propionyl	189	C12H15NO	1000123-86-3	10
4		(E)-3-Chloro-2-methyl-2-pentenal	132	C6H9ClO	031357-76-3	9
5		1,3,5-Trifluorobenzene	132	C6H3F3	000372-38-3	9



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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Butane, 2-methoxy...	2.19	2.3	ng	90160	1	6.90	796636	20.0
unknown6.67	6.67	36.2	ng	1440340	1	6.90	796636	20.0