

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
 Data File : BF088122.D
 Acq On : 18 Jun 2016 6:00
 Operator : UM/SJ
 Sample : H3574-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-26

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:\HPCHEM1\BNA_F\METHODS\8270-BF060716.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	1.678	6	8	16	rVV	99888	241472	7.67%	0.368%
2	1.793	16	18	39	rVB	1117820	2090924	66.44%	3.188%
3	2.513	73	81	84	rVB	30784	70687	2.25%	0.108%
4	2.684	93	96	98	rBV	72407	139303	4.43%	0.212%
5	2.718	98	99	105	rVB	102247	165109	5.25%	0.252%
6	3.816	191	195	201	rVB2	19954	46230	1.47%	0.070%
7	4.056	213	216	218	rBV	81994	130349	4.14%	0.199%
8	4.753	274	277	279	rBV2	32295	64301	2.04%	0.098%
9	4.856	284	286	288	rVV	45979	76700	2.44%	0.117%
10	4.901	288	290	294	rVB2	35813	65321	2.08%	0.100%
11	5.290	323	324	329	rBV	878248	1103351	35.06%	1.682%
12	5.404	332	334	337	rBV3	21982	43963	1.40%	0.067%
13	5.587	347	350	353	rVB	39483	47942	1.52%	0.073%
14	5.827	368	371	373	rBV	70737	93854	2.98%	0.143%
15	5.919	377	379	382	rVB	80151	74481	2.37%	0.114%
16	6.204	401	404	407	rBV3	27364	73901	2.35%	0.113%
17	6.296	410	412	413	rBV	38018	44335	1.41%	0.068%
18	6.330	413	415	416	rBV	51554	78328	2.49%	0.119%
19	6.353	416	417	421	rBV	506500	630315	20.03%	0.961%
20	6.433	421	424	425	rVB2	79835	146296	4.65%	0.223%
21	6.467	425	427	429	rBV	1656762	1806050	57.39%	2.754%
22	6.513	429	431	433	rVB	110171	97210	3.09%	0.148%
23	6.593	435	438	440	rVB2	47030	68211	2.17%	0.104%
24	6.650	440	443	445	rBV3	66408	84837	2.70%	0.129%
25	6.696	445	447	449	rBV	678546	605340	19.24%	0.923%
26	6.776	452	454	455	rVB	419976	370902	11.79%	0.566%
27	6.856	458	461	464	rBV2	1858511	2517773	80.01%	3.839%
28	6.947	466	469	471	rBV	460749	502657	15.97%	0.766%
29	6.982	471	472	474	rVB	43636	46864	1.49%	0.071%
30	7.039	474	477	480	rVB3	300737	395541	12.57%	0.603%
31	7.085	480	481	482	rBV	51345	49555	1.57%	0.076%
32	7.142	482	486	487	rBV2	92450	110684	3.52%	0.169%
33	7.245	489	495	496	rBV2	846539	1650418	52.45%	2.517%
34	7.267	496	497	499	rVB2	1294394	1587364	50.44%	2.420%

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35	7.347	499	504	506	rVB4	283714	569604	18.10%	0.869%
36	7.393	506	508	510	rVB3	287476	489295	15.55%	0.746%
37	7.473	513	515	517	rVB	764572	710358	22.57%	1.083%
38	7.530	517	520	522	rBV4	169746	364742	11.59%	0.556%
39	7.565	522	523	524	rVB	142294	97614	3.10%	0.149%
40	7.599	524	526	527	rVB2	58414	73927	2.35%	0.113%
41	7.645	527	530	533	rBV3	331783	677828	21.54%	1.034%
42	7.690	533	534	535	rBV	100732	92649	2.94%	0.141%
43	7.725	535	537	542	rVB	1547359	1978813	62.88%	3.017%
44	7.816	542	545	547	rBV3	180647	386495	12.28%	0.589%
45	7.862	547	549	551	rVB	177557	196370	6.24%	0.299%
46	7.930	551	555	556	rBV3	242383	535863	17.03%	0.817%
47	7.987	556	560	563	rBV2	1011196	2070314	65.79%	3.157%
48	8.033	563	564	565	rVB	469528	322096	10.24%	0.491%
49	8.079	567	568	572	rVB3	223001	410233	13.04%	0.626%
50	8.148	572	574	575	rBV	134803	206818	6.57%	0.315%
51	8.182	575	577	580	rBV3	336149	586063	18.62%	0.894%
52	8.262	580	584	587	rBV5	254419	871565	27.70%	1.329%
53	8.308	587	588	590	rVB2	198884	229977	7.31%	0.351%
54	8.365	592	593	595	rVB	220113	222485	7.07%	0.339%
55	8.422	595	598	599	rBV3	305534	498165	15.83%	0.760%
56	8.456	599	601	602	rVV2	330011	422065	13.41%	0.644%
57	8.490	602	604	606	rVV2	422585	549402	17.46%	0.838%
58	8.525	606	607	608	rVB	222821	155526	4.94%	0.237%
59	8.616	613	615	616	rBV2	239356	346259	11.00%	0.528%
60	8.650	616	618	620	rVB2	337527	444056	14.11%	0.677%
61	8.696	620	622	623	rBV2	294920	311370	9.89%	0.475%
62	8.765	626	628	629	rBV	229074	288509	9.17%	0.440%
63	8.799	629	631	633	rBV	1770099	1968724	62.56%	3.002%
64	8.845	633	635	636	rVB2	224926	256299	8.14%	0.391%
65	8.879	636	638	640	rBV2	181793	297659	9.46%	0.454%
66	9.028	649	651	652	rBV2	378244	579747	18.42%	0.884%
67	9.062	652	654	656	rVV	1768502	2845606	90.42%	4.339%
68	9.119	656	659	661	rVB4	258478	510283	16.22%	0.778%
69	9.165	661	663	664	rBV	237191	238391	7.58%	0.364%
70	9.256	664	671	674	rBV4	599899	1895182	60.22%	2.890%
71	9.325	674	677	681	rVB	996087	1798179	57.14%	2.742%

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72	9.405	681	684	688	rBV2	1405798	3146942	100.00%	4.799%
73	9.519	691	694	698	rVB2	1024257	2262099	71.88%	3.449%
74	9.599	698	701	704	rVB5	490942	833812	26.50%	1.271%
75	9.645	704	705	707	rBV2	326259	424633	13.49%	0.647%
76	9.713	709	711	712	rBV	222721	369404	11.74%	0.563%
77	9.736	712	713	714	rBV	642555	588571	18.70%	0.897%
78	9.839	720	722	724	rBV3	439874	529794	16.84%	0.808%
79	9.885	724	726	728	rBV2	247958	506623	16.10%	0.773%
80	9.931	728	730	733	rVV2	684931	1298052	41.25%	1.979%
81	9.976	733	734	738	rVB3	243093	568756	18.07%	0.867%
82	10.045	738	740	742	rVB2	351526	631891	20.08%	0.964%
83	10.091	742	744	746	rBV2	474263	877823	27.89%	1.339%
84	10.148	746	749	750	rVV	674800	1113678	35.39%	1.698%
85	10.182	750	752	755	rVB3	576184	963821	30.63%	1.470%
86	10.239	755	757	759	rBV3	317312	590605	18.77%	0.901%
87	10.365	766	768	773	rBV5	474515	1131685	35.96%	1.726%
88	10.445	773	775	777	rVV	393646	561369	17.84%	0.856%
89	10.536	781	783	784	rVV2	882262	987435	31.38%	1.506%
90	10.571	784	786	789	rVV3	653924	1295034	41.15%	1.975%
91	10.616	789	790	795	rVV2	506464	1086266	34.52%	1.656%
92	10.719	797	799	801	rBV2	249842	368931	11.72%	0.563%
93	10.856	810	811	813	rVB	440911	456480	14.51%	0.696%
94	10.959	819	820	824	rBV3	254624	523865	16.65%	0.799%
95	11.108	831	833	838	rVV4	418488	726110	23.07%	1.107%
96	11.222	841	843	846	rVB	826714	840186	26.70%	1.281%
97	12.148	922	924	926	rBV	175650	183903	5.84%	0.280%
98	12.799	979	981	984	rVB	1394359	1659332	52.73%	2.530%
99	13.840	1070	1072	1075	rVB	372129	523964	16.65%	0.799%
100	15.223	1191	1193	1209	rVB	405710	710630	22.58%	1.084%

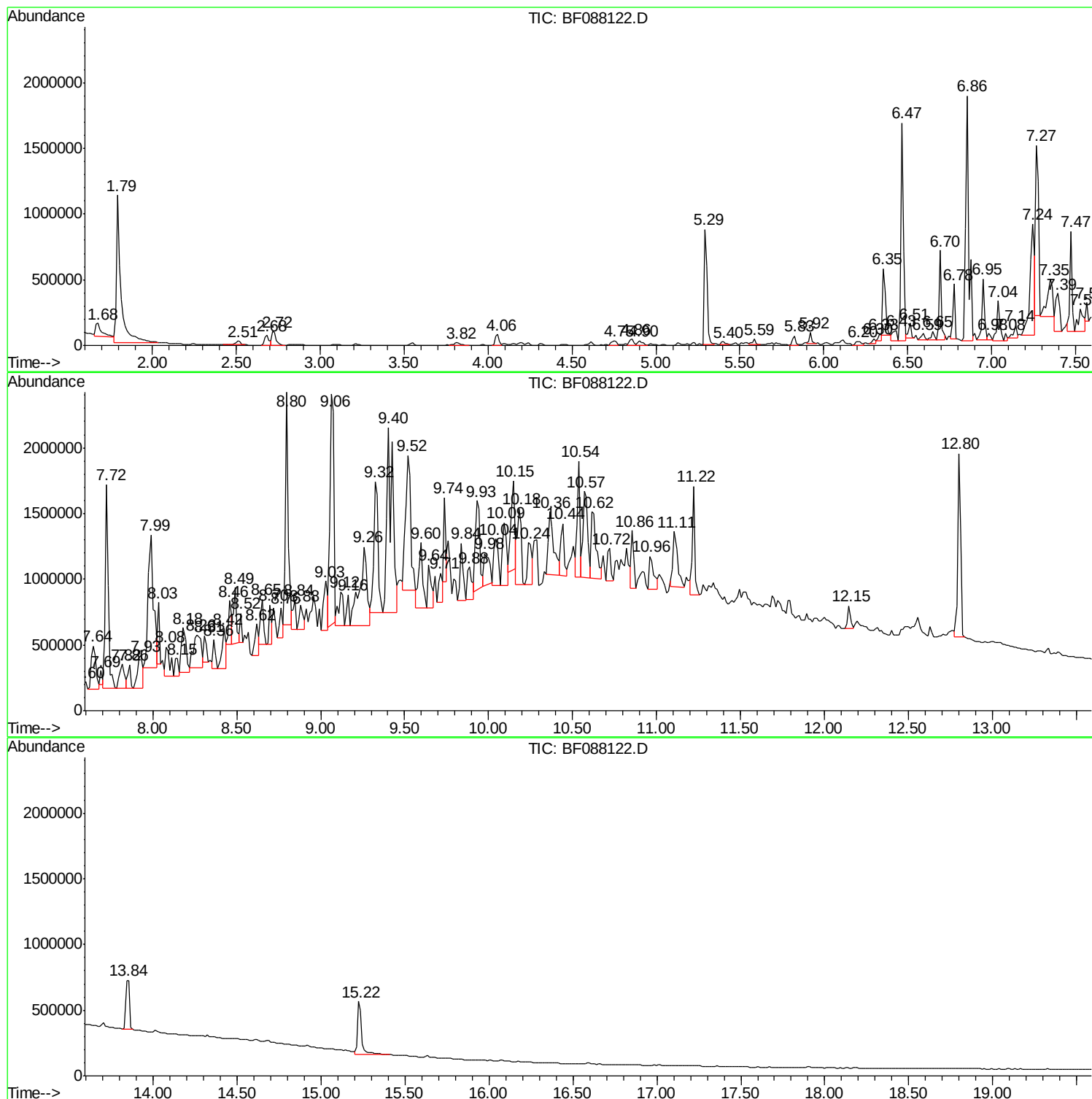
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MW-26

Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P



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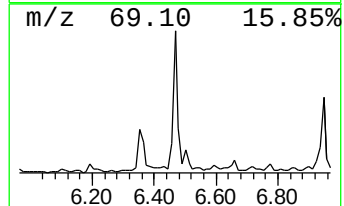
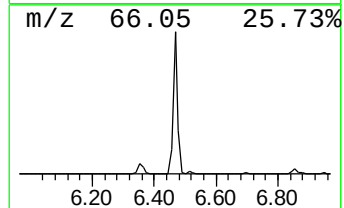
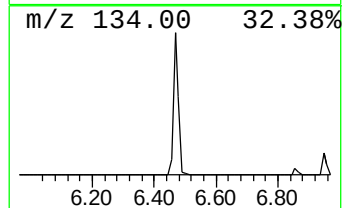
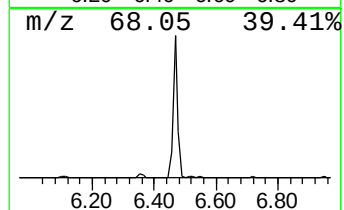
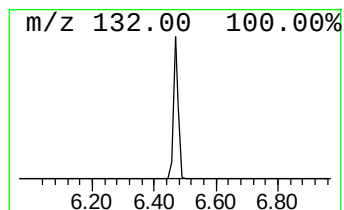
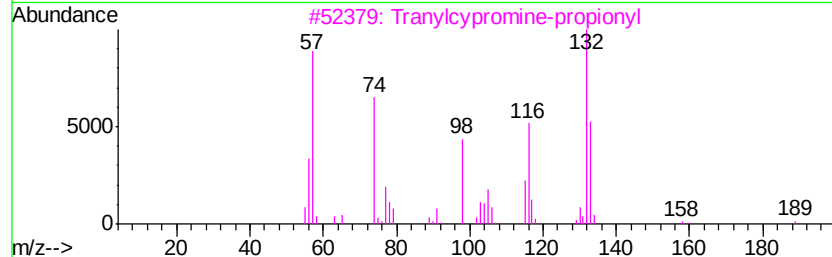
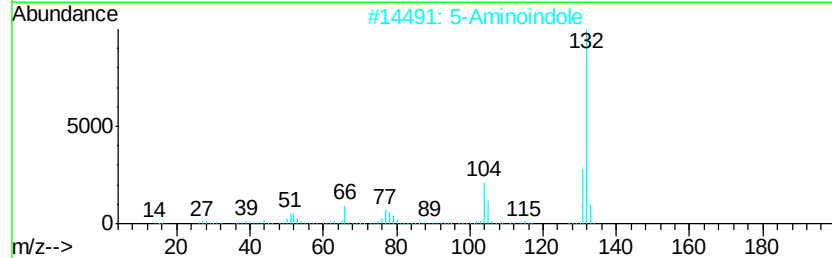
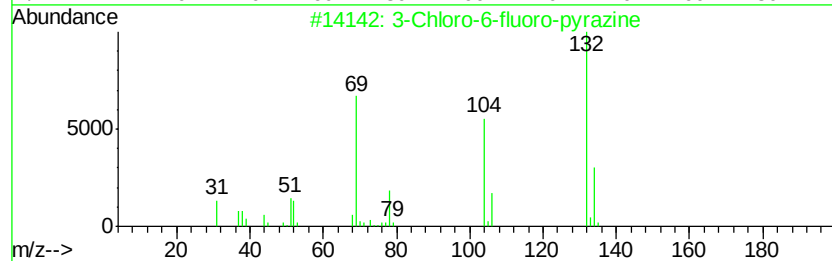
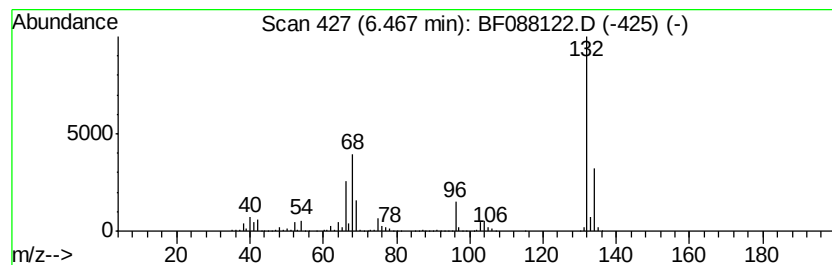
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.47 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.47	59.67 ng	1806050	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Chloro-6-fluoro-pyrazine			132	C4H2ClFN2	1000146-10-7	27
2	5-Aminoindole			132	C8H8N2	005192-03-0	12
3	Tranvlcyproline-propionyl			189	C12H15NO	1000123-86-3	10
4	Bicyclo[4.2.0]octa-1,3,5-triene,...			132	C10H12	028749-81-7	9
5	(E)-3-Chloro-2-methyl-2-pentenal			132	C6H9ClO	031357-76-3	9



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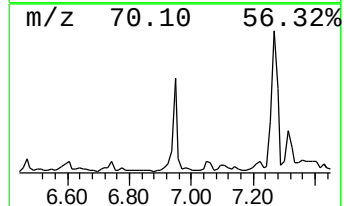
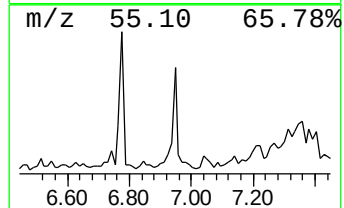
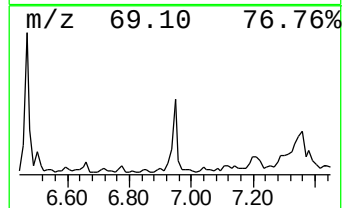
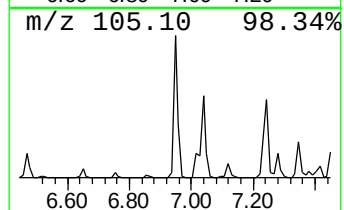
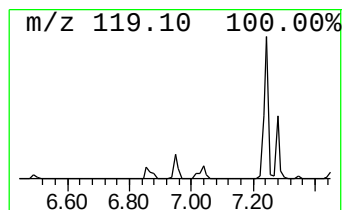
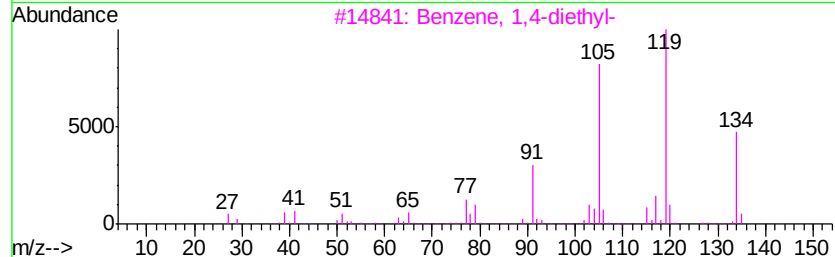
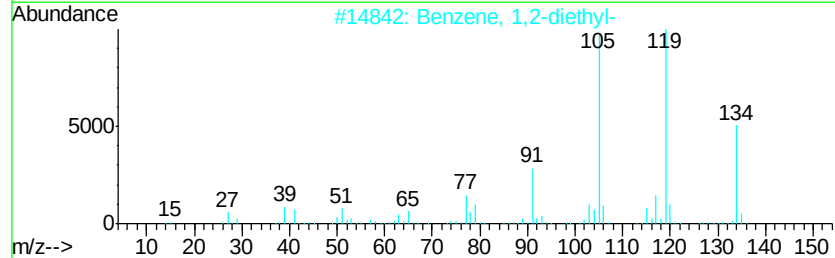
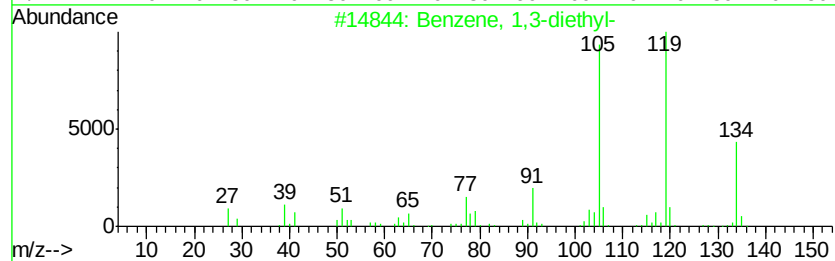
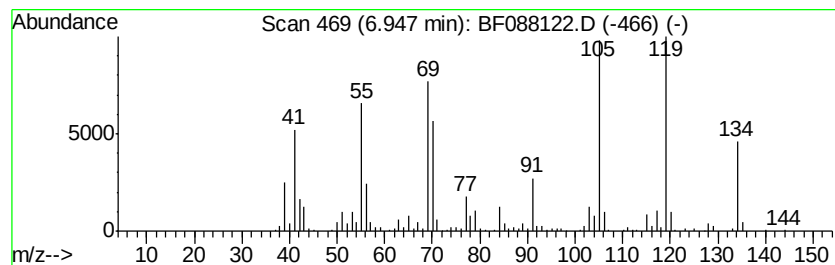
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TIC Library : C:\Database\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, 1,3-diethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.95	16.61 ng	502657	1,4-Dichlorobenzene-d4	6.70

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	95
3			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	90
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	55
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	50



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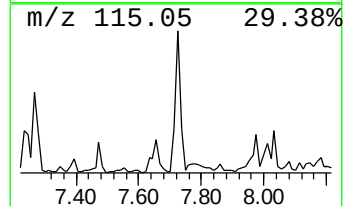
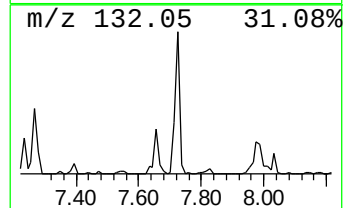
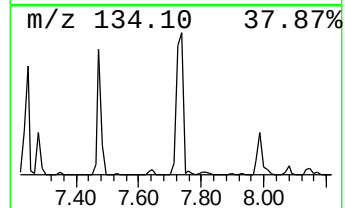
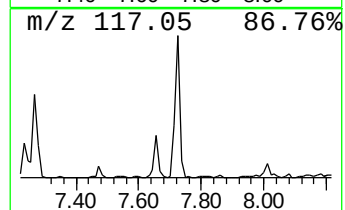
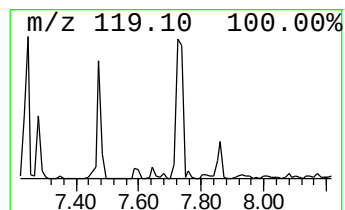
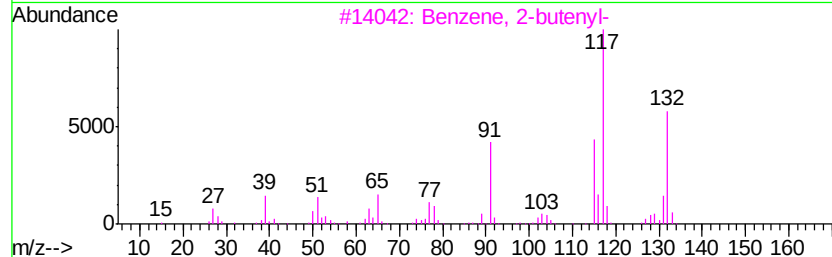
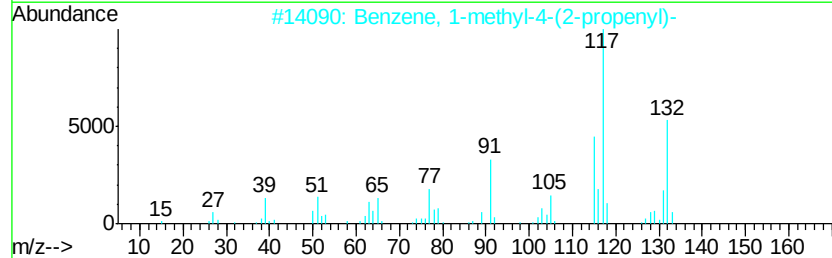
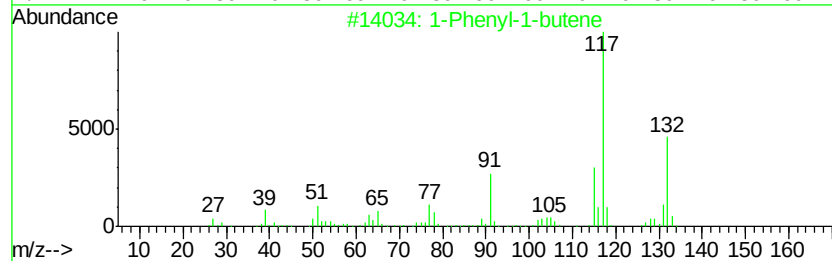
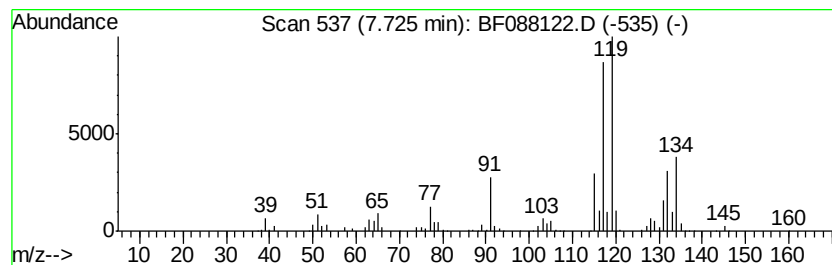
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Peak Number 5 1-Phenyl-1-butene Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.72	19.12 ng	1978810	Naphthalene-d8	7.99

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Phenyl-1-butene	132	C10H12	000824-90-8	91
2		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3		Benzene, 2-butenyl-	132	C10H12	001560-06-1	70
4		Benzene, (1-methyl-1-propenyl)-, ...	132	C10H12	000768-00-3	70
5		Benzene, 1-ethenyl-3-ethyl-	132	C10H12	007525-62-4	55



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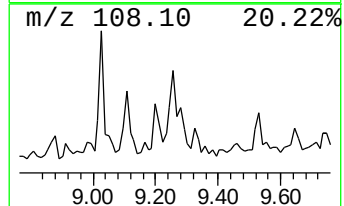
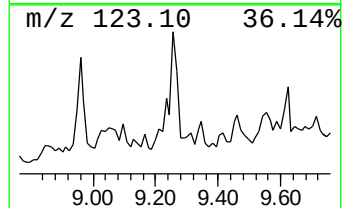
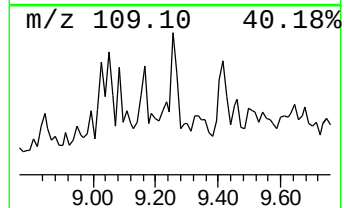
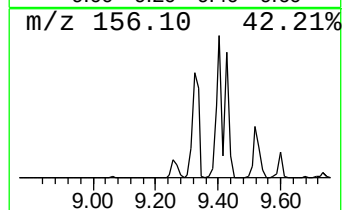
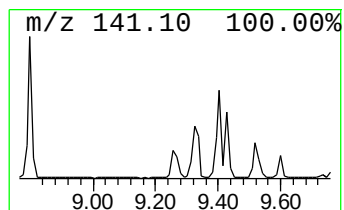
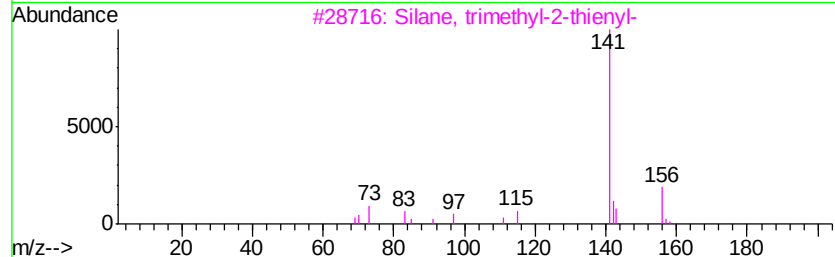
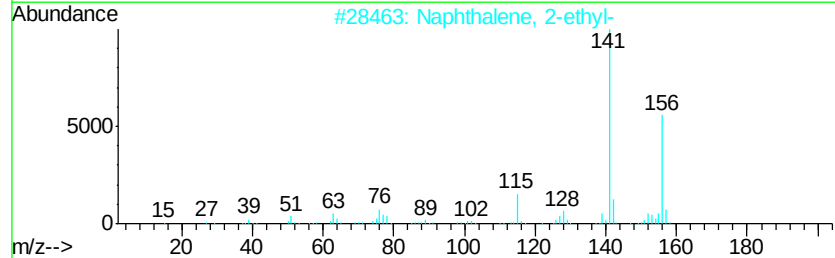
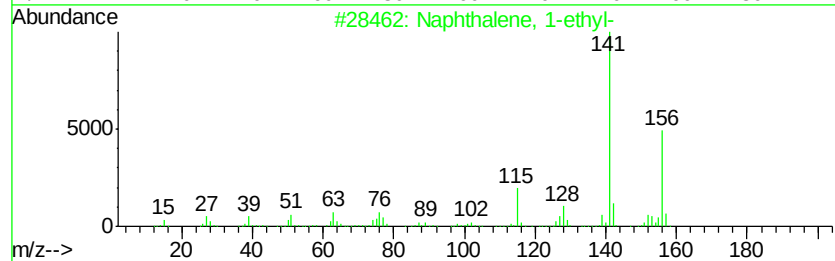
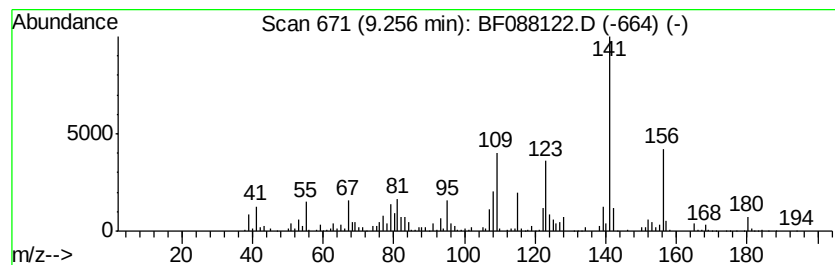
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Naphthalene, 1-ethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.26	64.40 ng	1895180	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	60
2		Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	60
3		Silane, trimethyl-2-thienyl-	156	C7H12SSi	018245-28-8	43
4		Naphthalene, 1-(chloromethyl)-	176	C11H9Cl	000086-52-2	38
5		2,5-Dimethoxyfluorobenzene	156	C8H9FO2	082830-49-7	38



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-26

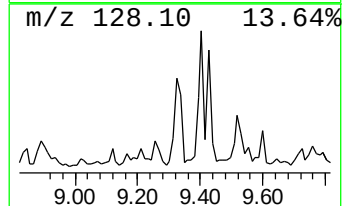
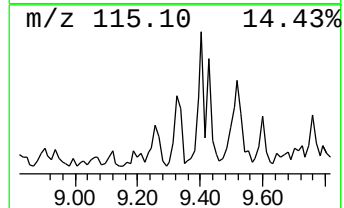
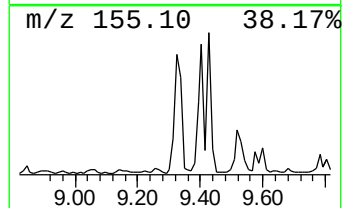
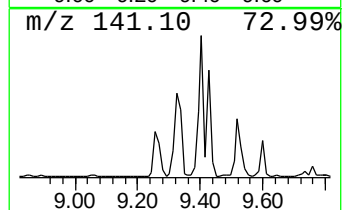
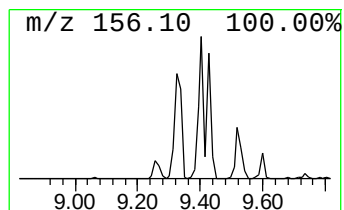
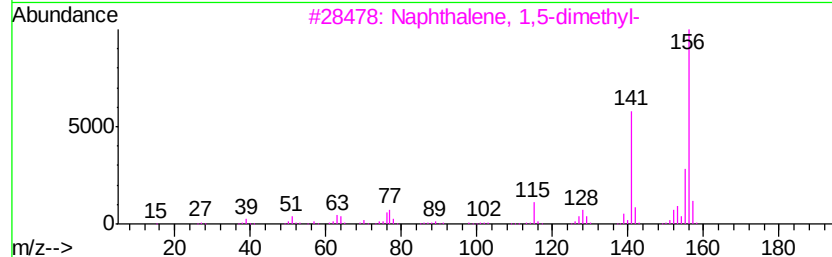
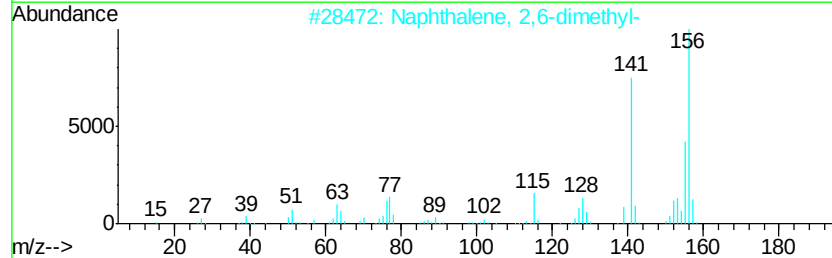
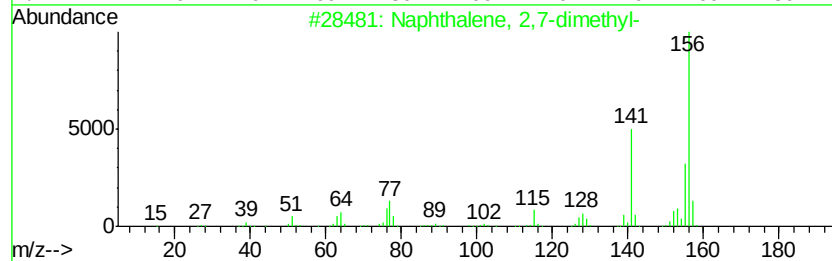
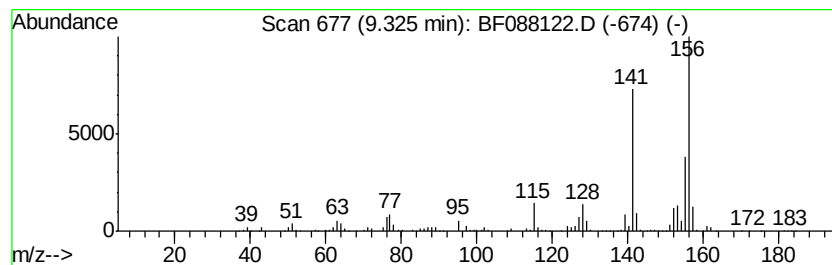
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Naphthalene, 2,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.32	61.10 ng	1798180	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97
3		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
4		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5		Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-26

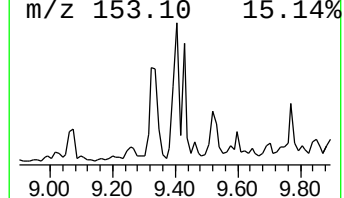
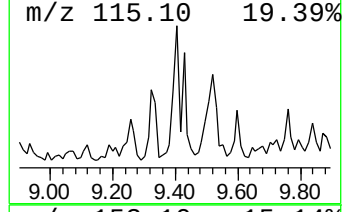
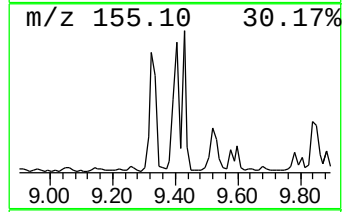
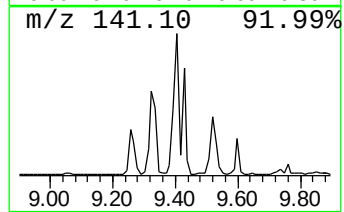
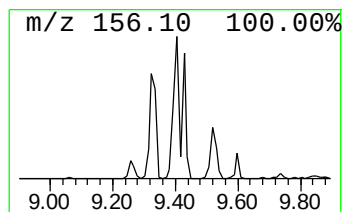
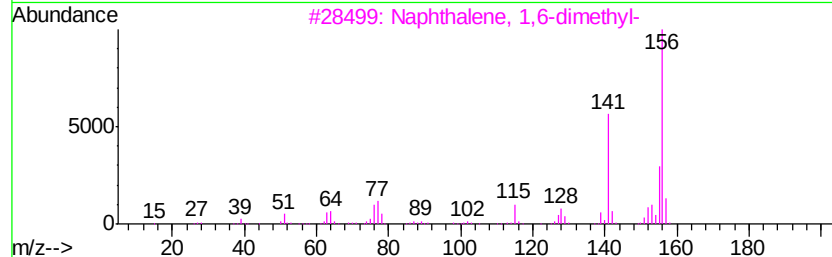
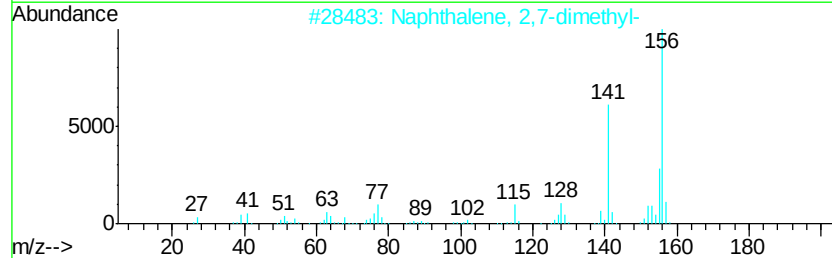
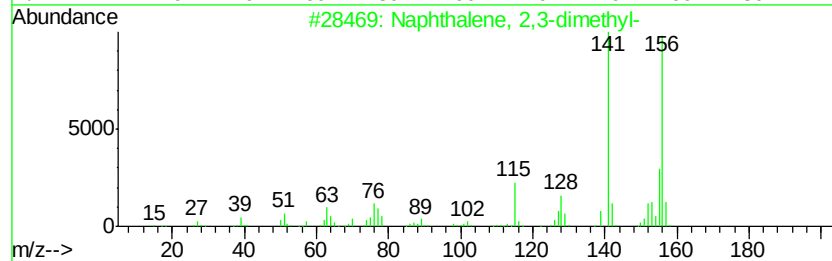
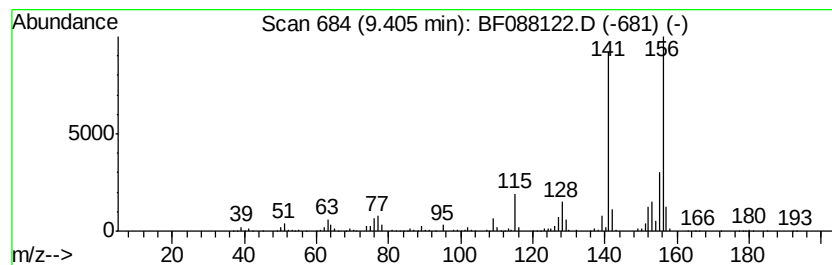
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Naphthalene, 2,3-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.40	106.94 ng	3146940	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2		Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3		Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
4		Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	97
5		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	97



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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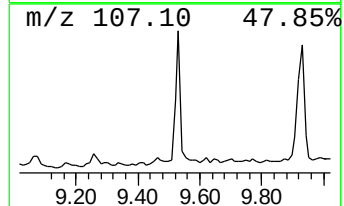
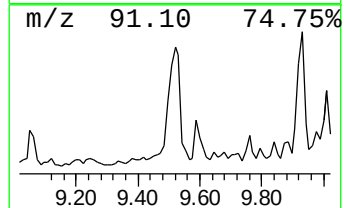
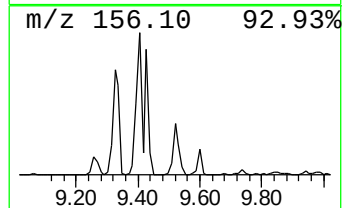
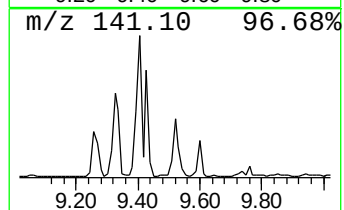
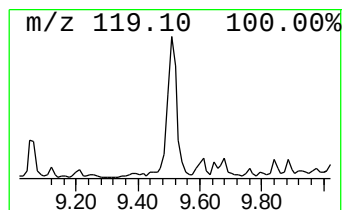
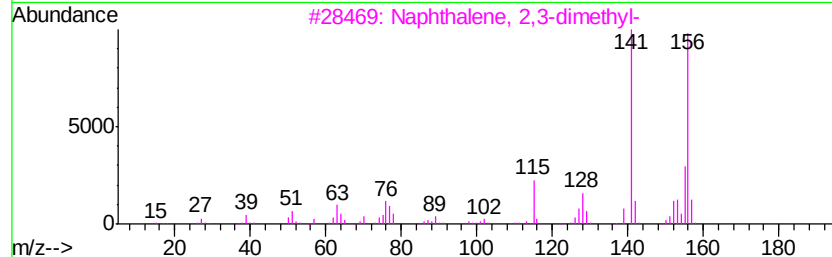
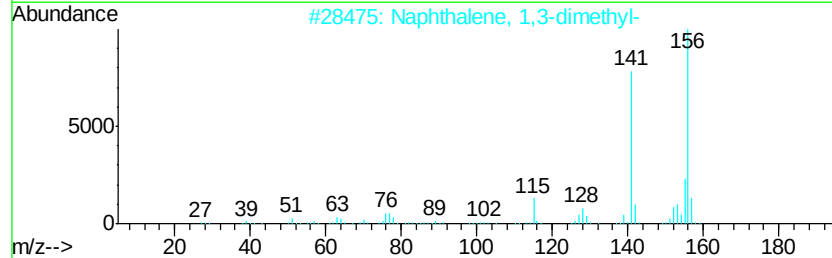
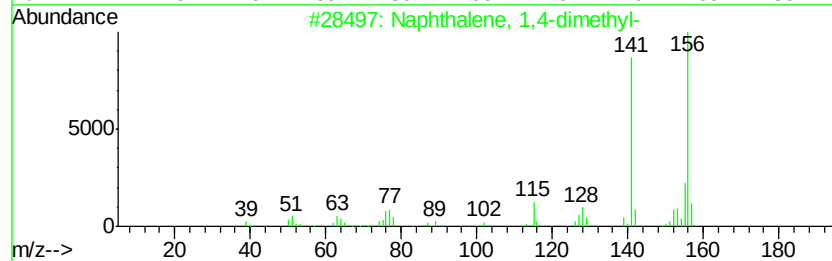
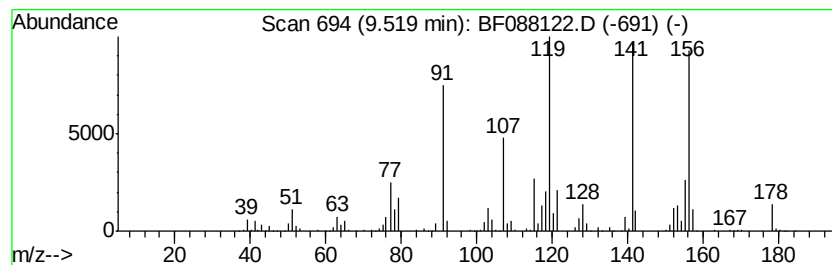
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Naphthalene, 1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.52	76.87 ng	2262100	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2		Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
3		Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	90
4		Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	86
5		Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	86



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampled :
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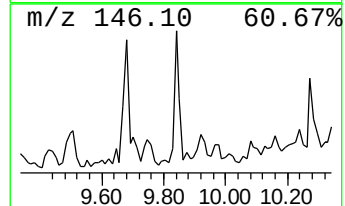
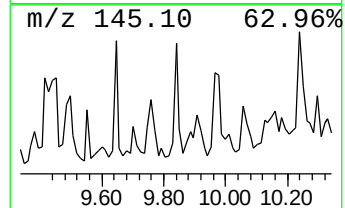
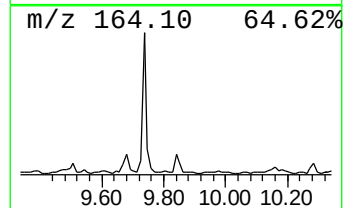
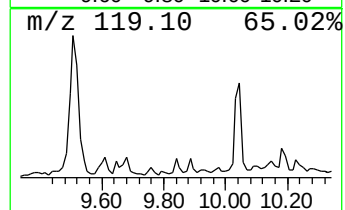
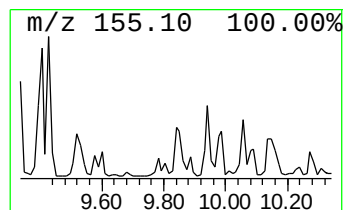
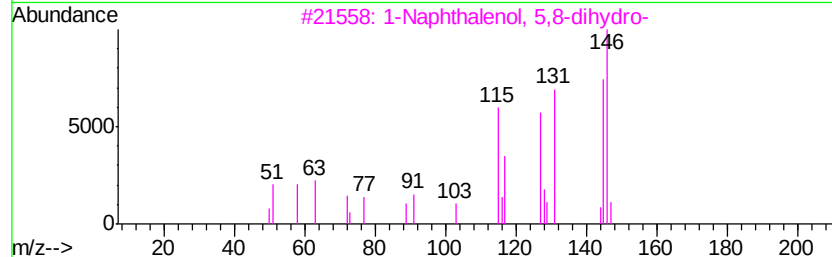
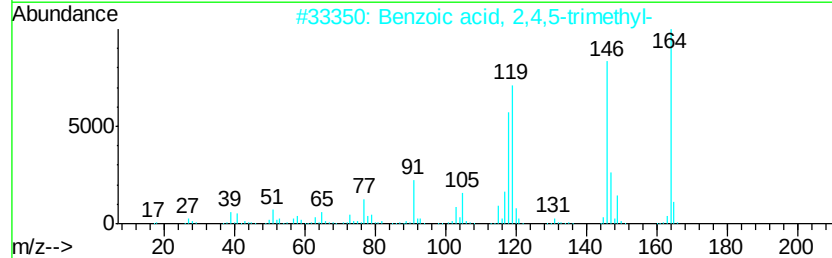
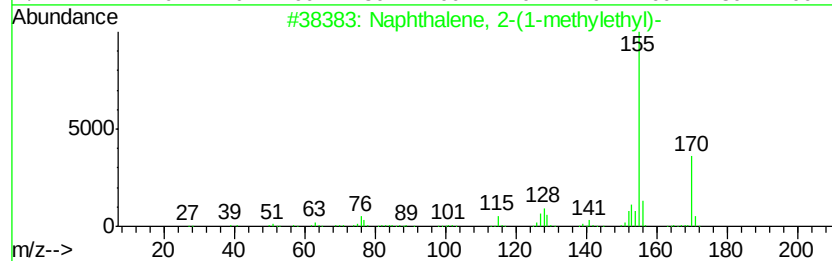
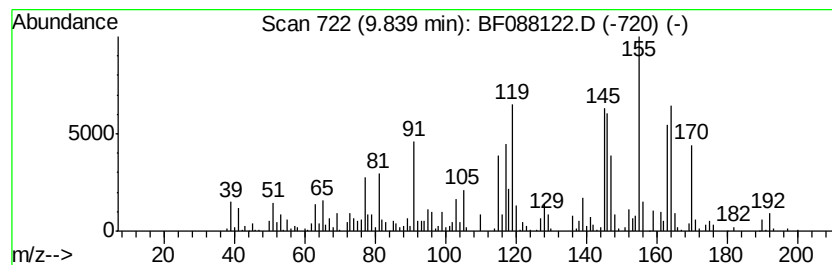
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown9.84 Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.84	18.00 ng	529794	Acenaphthene-d10	9.74

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	35
2			Benzoic acid, 2,4,5-trimethyl-	164	C10H12O2	000528-90-5	30
3			1-Naphthalenol, 5,8-dihydro-	146	C10H10O	027673-48-9	25
4			2-Propenoic acid, 3-(3-hydroxyph...	164	C9H8O3	000588-30-7	20
5			2-Propenal, 2-methyl-3-phenyl-	146	C10H10O	000101-39-3	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
MW-26

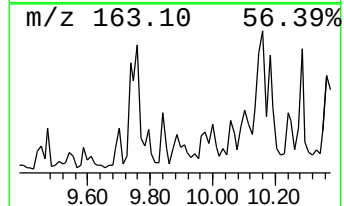
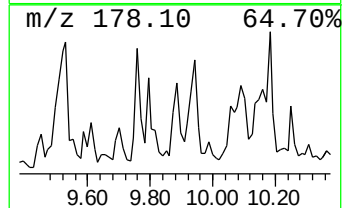
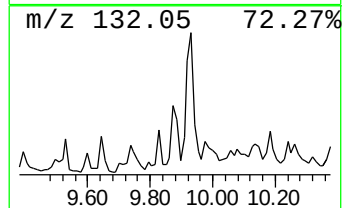
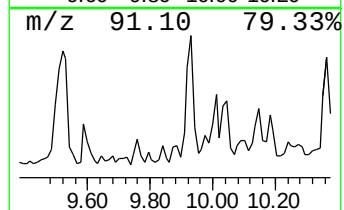
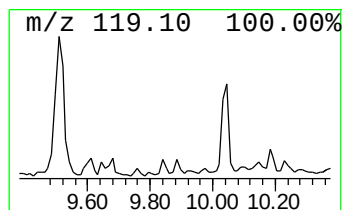
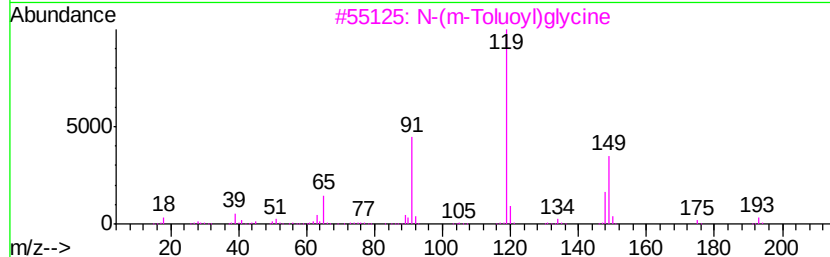
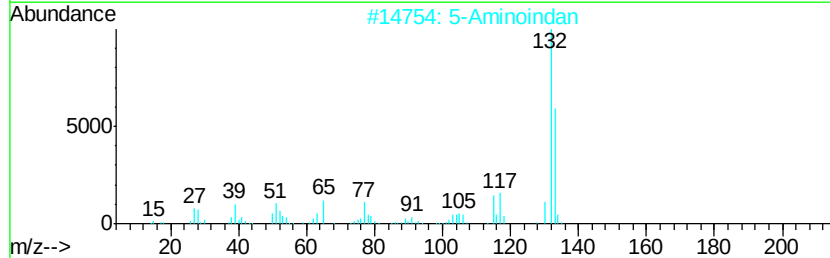
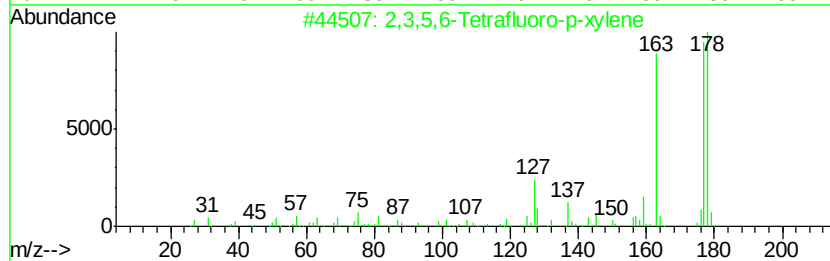
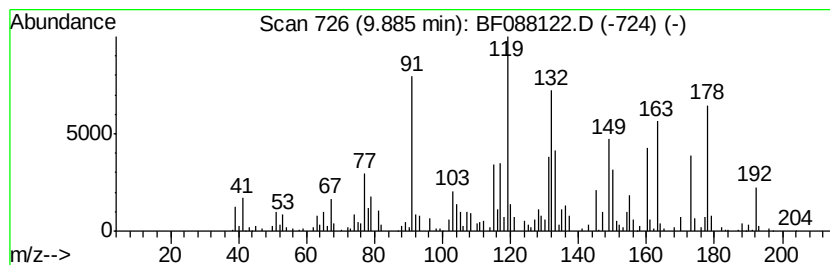
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 unknown9.88 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.88	17.22 ng	506623	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2,3,5,6-Tetrafluoro-p-xylene	178	C8H6F4	000703-87-7	25
2		5-Aminoindan	133	C9H11N	024425-40-9	25
3		N-(m-Toluoyl)glycine	193	C10H11NO3	027115-49-7	18
4		6-Isobutyl-4-methyl-5,6-dihydro-...	176	C11H16N2	1000185-72-5	18
5		Benzeneacetic acid, .alpha.-ethy...	178	C11H14O2	002294-71-5	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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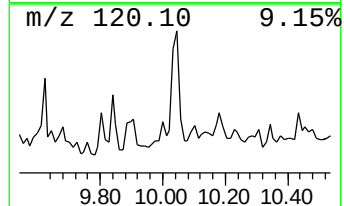
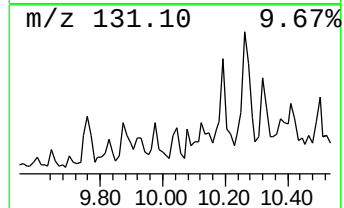
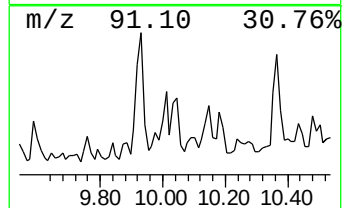
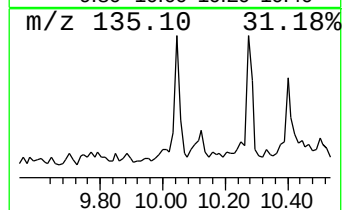
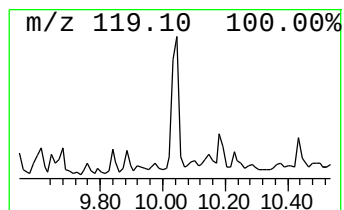
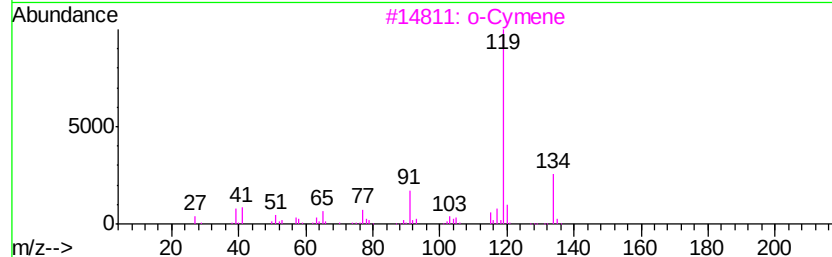
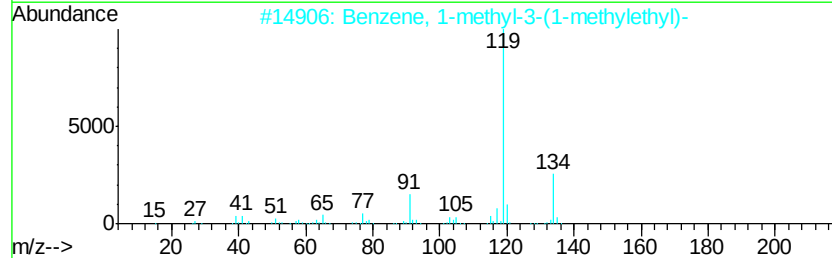
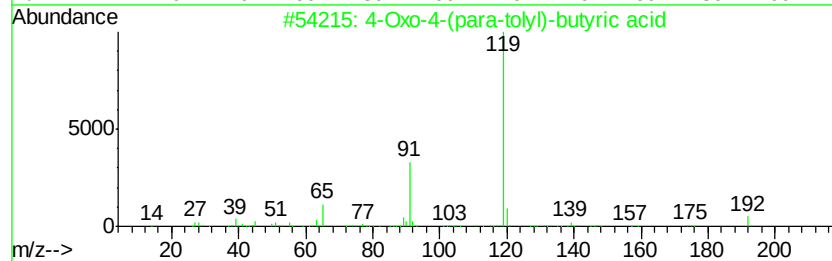
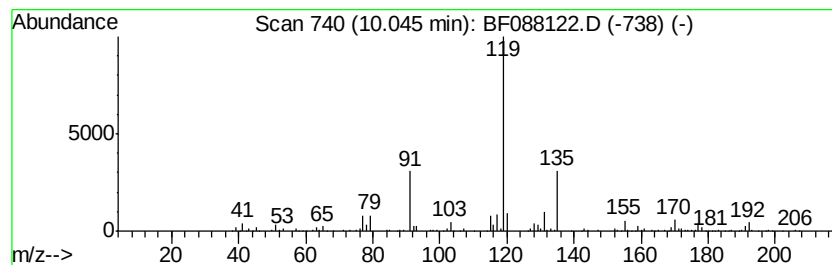
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 4-Oxo-4-(para-tolyl)-butyri... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.04	21.47 ng	631891	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	4-Oxo-4-(para-tolyl)-butyric acid	192	C11H12O3	004619-20-9	50
2		Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	49
3		o-Cymene	134	C10H14	000527-84-4	47
4		Benzoic acid, 3-methyl-, 2,4,5-trimethyl-	314	C14H9Cl3O2	1000355-71-4	47
5		2-Methylbenzoic acid, 2-fluorophenyl-	230	C14H11FO2	1000325-60-5	47



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-26

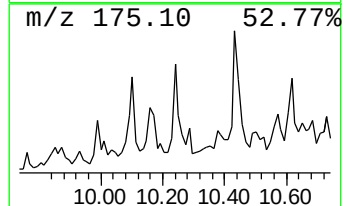
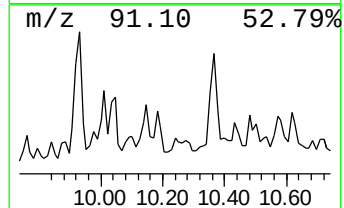
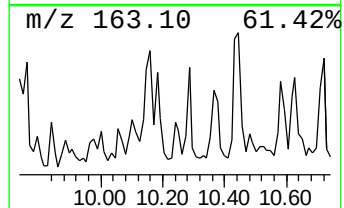
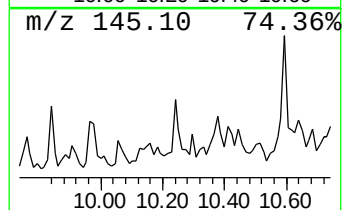
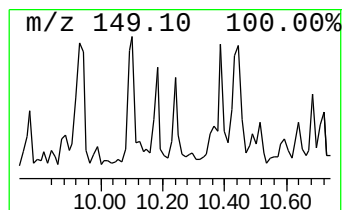
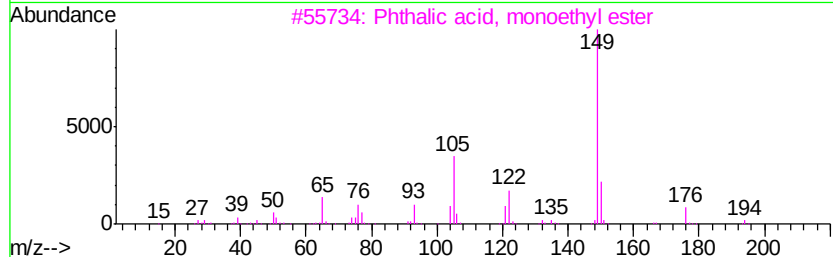
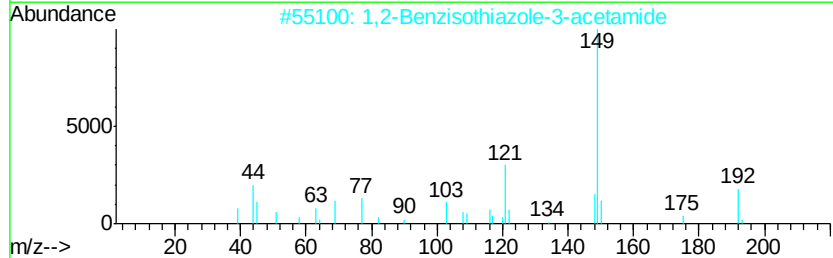
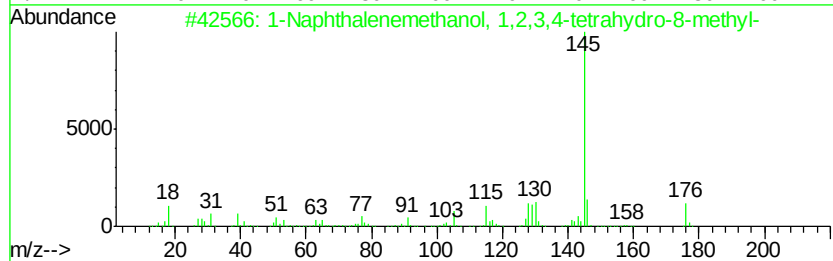
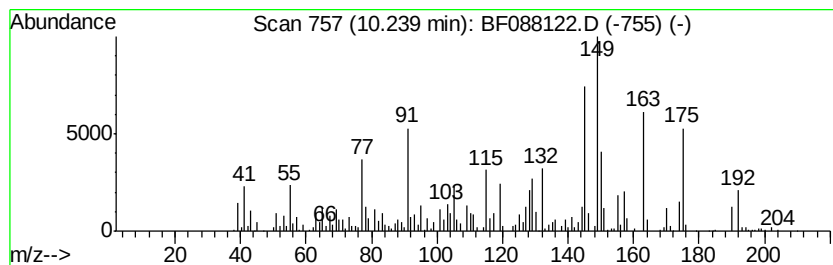
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 16 unknown10.24 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.24	20.07 ng	590605	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1-Naphthalenemethanol, 1,2,3,4-t...	176	C12H16O	036052-28-5	25
2		1,2-Benzisothiazole-3-acetamide	192	C9H8N2OS	029273-65-2	14
3		Phthalic acid, monoethyl ester	194	C10H10O4	002306-33-4	11
4		1-Phenylcyclopentanecarboxylic acid	190	C12H14O2	000077-55-4	11
5		Quinazolin-6-ylamine	145	C8H7N3	101421-72-1	10



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-26

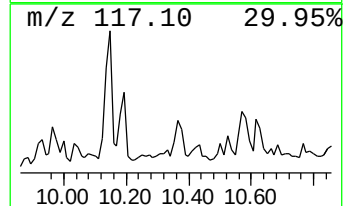
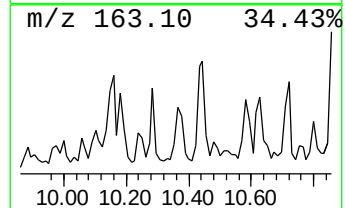
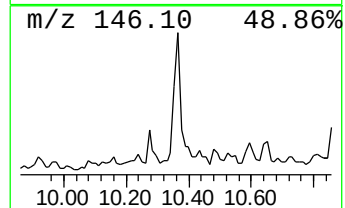
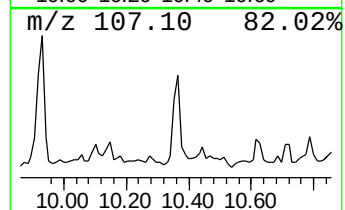
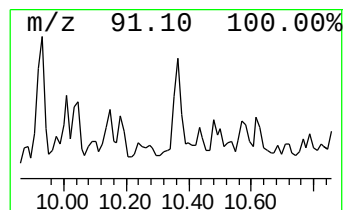
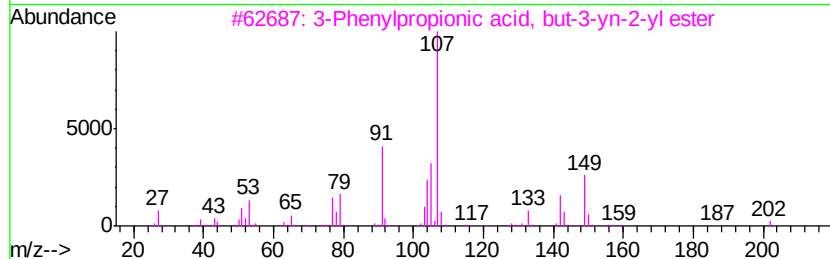
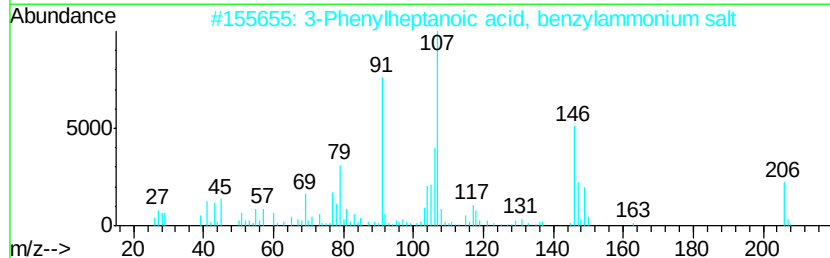
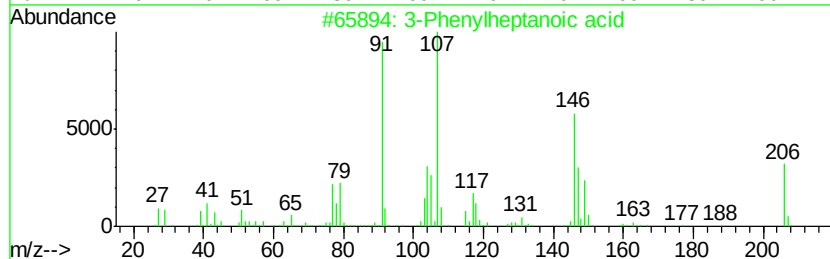
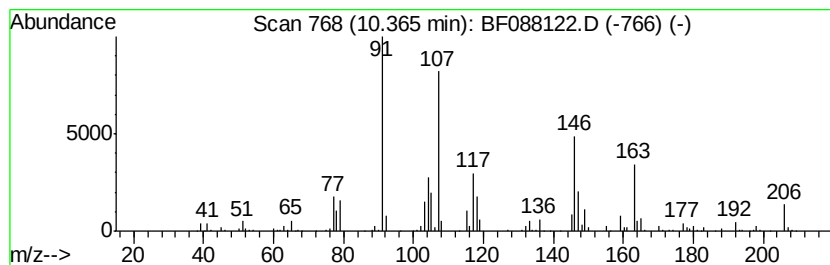
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 3-Phenylheptanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.36	38.46 ng	1131690	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	3-Phenylheptanoic acid	206	C13H18O2	005638-30-2	62
2		3-Phenylheptanoic acid, benzylam...	313	C20H27NO2	1000196-52-4	50
3		3-Phenylpropionic acid, but-3-yn...	202	C13H14O2	1000299-17-2	27
4		p-Propionotoluidide	163	C10H13NO	002759-55-9	27
5		Benzenemethanol, .alpha.-(bromom...	200	C8H9BrO	002425-28-7	27



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
MW-26

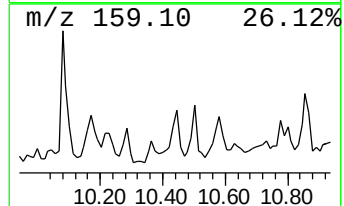
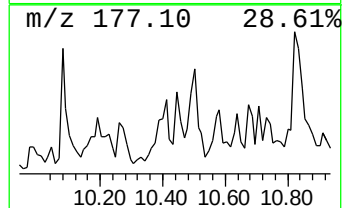
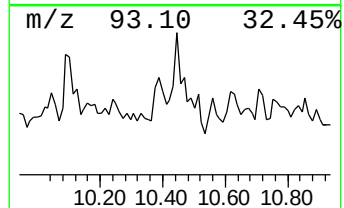
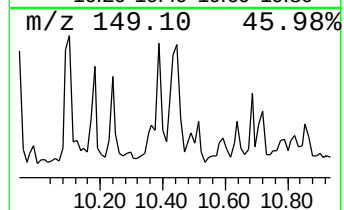
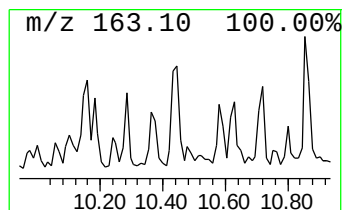
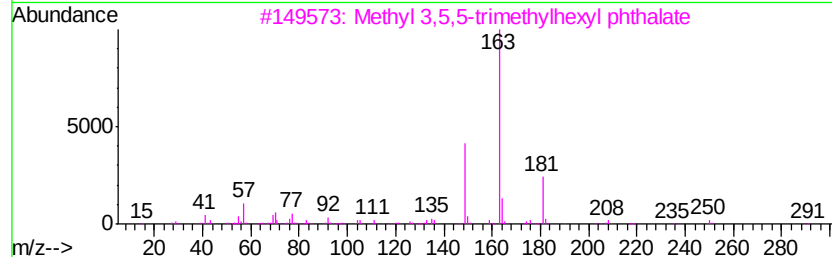
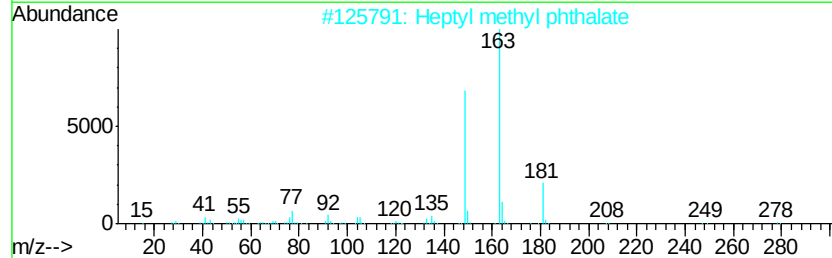
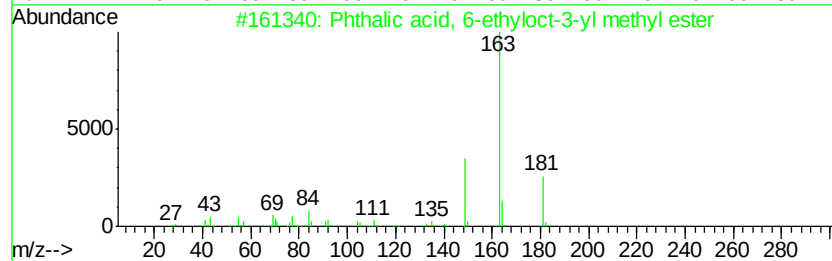
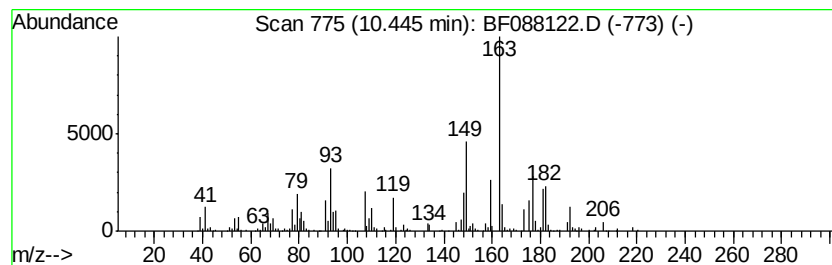
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 18 unknown10.44 Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.44	19.08 ng	561369	Acenaphthene-d10	9.74

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Phthalic acid, 6-ethyloct-3-yl m...	320	C19H28O4	1000315-53-0	43
2		Heptyl methyl phthalate	278	C16H22O4	1000373-88-4	43
3		Methyl 3,5,5-trimethylhexyl phth...	306	C18H26O4	1000373-81-4	43
4		Methyl 4-methylpentyl phthalate	264	C15H20O4	1000373-82-6	43
5		Phthalic acid, 5-methoxy-3-methy...	294	C16H22O5	1000314-88-2	43



Data Path : Z:\HPCHEM1\BNA F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

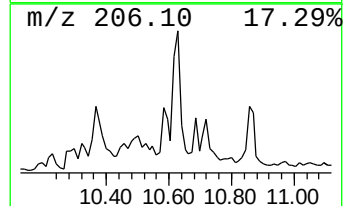
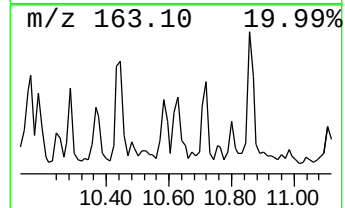
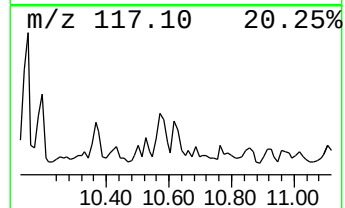
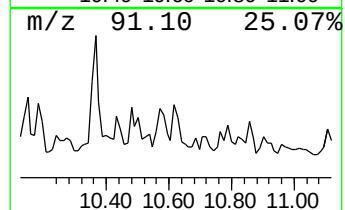
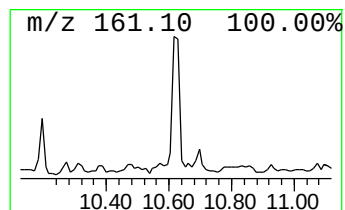
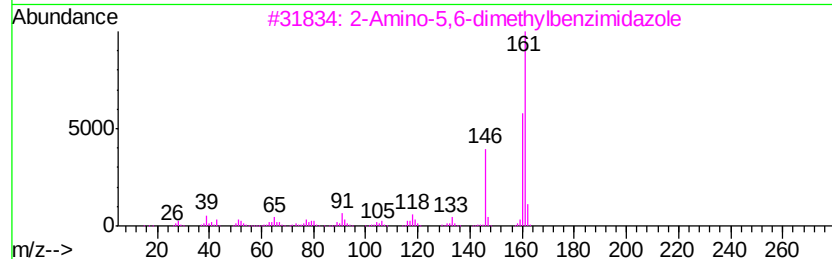
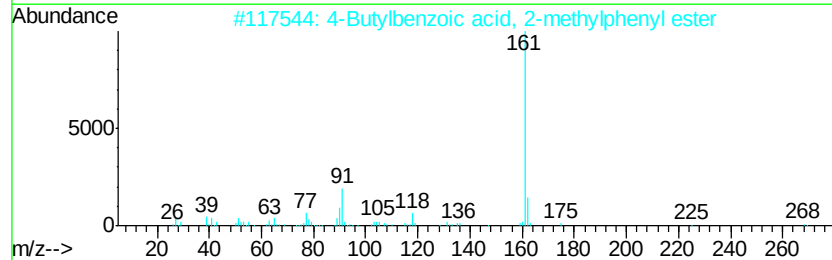
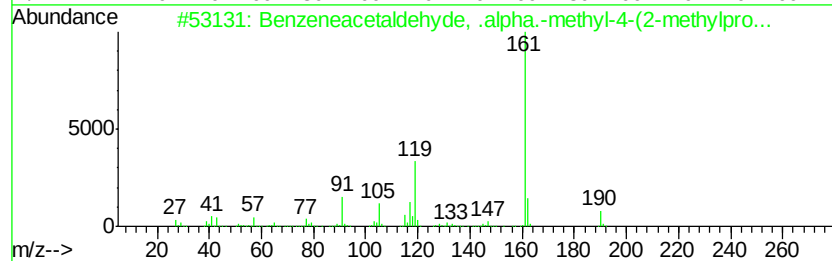
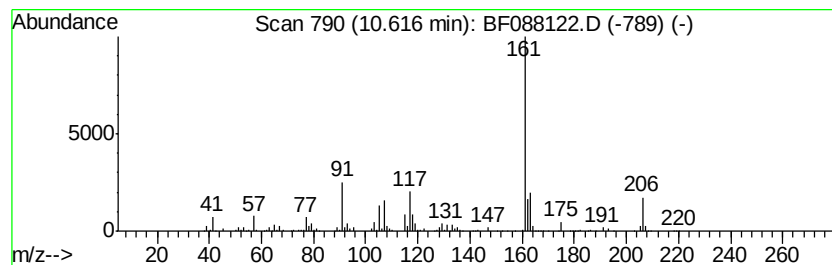
Instrument :
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ClientSampleId :
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Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 unknown10.62 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.		
10.62	25.86 ng	1086270	Phenanthrene-d10	11.22		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzeneacetaldehyde, .alpha.-met...	190	C13H18O	051407-46-6	47
2		4-Butylbenzoic acid, 2-methylphe...	268	C18H20O2	1000293-58-3	47
3		2-Amino-5,6-dimethylbenzimidazole	161	C9H11N3	029096-75-1	47
4		(7-Methylbenzo(b)thien-3-yl)acet...	206	C11H10O2S	001606-17-3	47
5		3-Isobutyl-4,5-dimethyl-3H-isobe...	218	C14H18O2	1000187-23-3	43



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
Data File : BF088122.D
Acq On : 18 Jun 2016 6:00
Operator : UM/SJ
Sample : H3574-06
Misc :
ALS Vial : 29 Sample Multiplier: 1

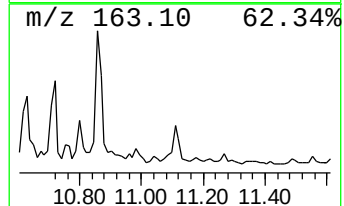
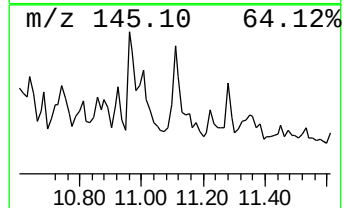
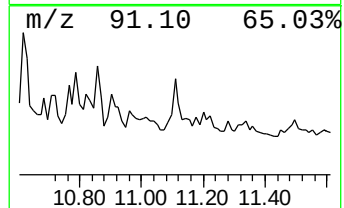
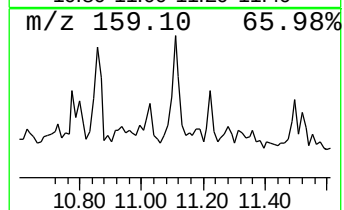
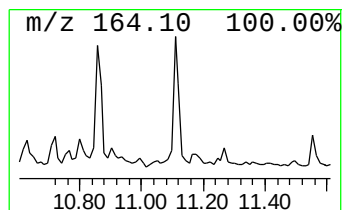
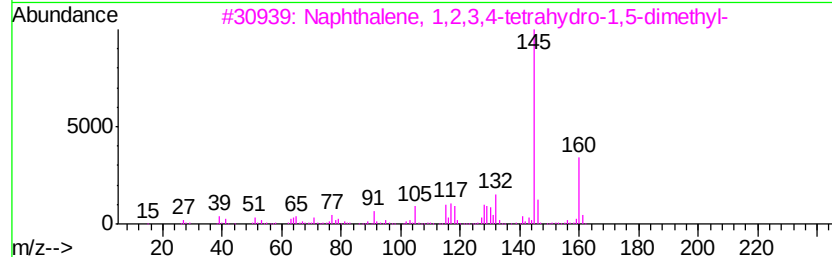
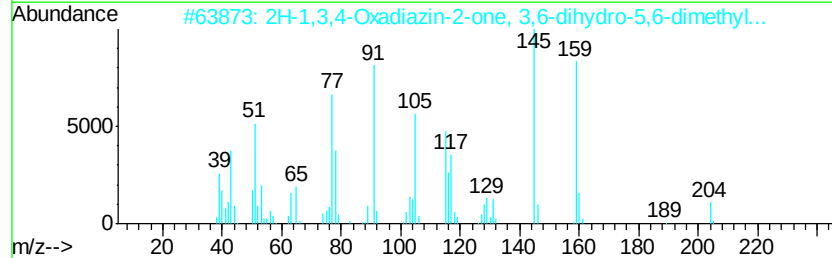
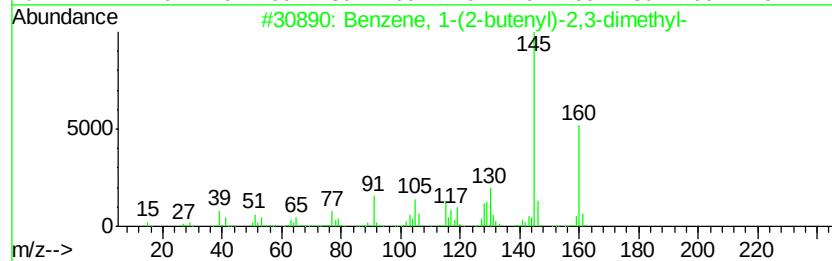
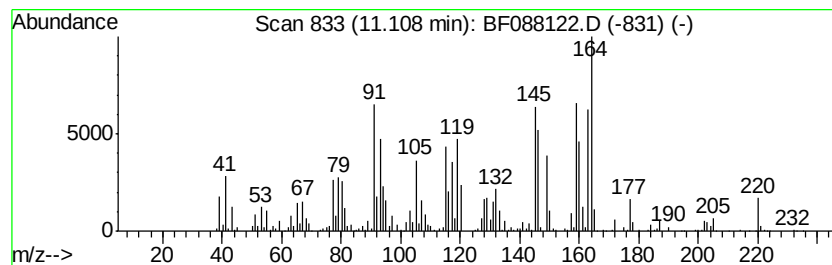
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Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 20 unknown11.11 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.11	17.28 ng	726110	Phenanthrene-d10	11.22	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	25
2	2H-1,3,4-Oxadiazin-2-one, 3,6-di...	204	C11H12N2O2	1000338-30-1	25
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	25
4	2-(5-Methyl-3-pyrrolyl)piperidine	164	C10H16N2	1000245-18-2	18
5	4-Methoxy-2-allylphenol	164	C10H12O2	1000306-52-7	18



Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061716\
 Data File : BF088122.D
 Acq On : 18 Jun 2016 6:00
 Operator : UM/SJ
 Sample : H3574-06
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 MW-26

Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF060716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Benzene, 1,3-diet...	6.95	16.6	ng	502657	1	6.70	605340	20.0
1-Phenyl-1-butene	7.72	19.1	ng	1978810	2	7.99	2070310	20.0
Naphthalene, 1-et...	9.26	64.4	ng	1895180	3	9.74	588571	20.0
Naphthalene, 2,7-...	9.32	61.1	ng	1798180	3	9.74	588571	20.0
Naphthalene, 2,3-...	9.40	106.9	ng	3146940	3	9.74	588571	20.0
Naphthalene, 1,4-...	9.52	76.9	ng	2262100	3	9.74	588571	20.0
unknown9.84	9.84	18.0	ng	529794	3	9.74	588571	20.0
unknown9.88	9.88	17.2	ng	506623	3	9.74	588571	20.0
4-Oxo-4-(para-tol...	10.04	21.5	ng	631891	3	9.74	588571	20.0
unknown10.24	10.24	20.1	ng	590605	3	9.74	588571	20.0
3-Phenylheptanoic...	10.36	38.5	ng	1131690	3	9.74	588571	20.0
unknown10.44	10.44	19.1	ng	561369	3	9.74	588571	20.0
unknown10.62	10.62	25.9	ng	1086270	4	11.22	840186	20.0
unknown11.11	11.11	17.3	ng	726110	4	11.22	840186	20.0