

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
 Data File : BF125023.D
 Acq On : 10 Aug 2021 2:59
 Operator : JU/CG
 Sample : M3314-01 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 72-11931

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-BF080421.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125023.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.022	497	499	502	rBB	2149	1712	4.13%	0.302%
2	5.434	566	569	573	rBB	46472	41459	100.00%	7.302%
3	6.451	739	742	747	rBB	47859	41461	100.00%	7.302%
4	6.822	802	805	808	rBB	38438	29361	70.82%	5.171%
5	7.381	897	900	903	rBB	34707	25494	61.49%	4.490%
6	8.004	1004	1006	1008	rBB	2035	1515	3.65%	0.267%
7	8.104	1020	1023	1026	rBB	45504	34078	82.19%	6.002%
8	9.175	1202	1205	1208	rBV	55538	40622	97.98%	7.154%
9	9.857	1318	1321	1324	rBB	44721	32409	78.17%	5.708%
10	10.645	1452	1455	1457	rBB	25755	18259	44.04%	3.216%
11	11.345	1571	1574	1576	rBV	37793	31914	76.97%	5.621%
12	11.363	1576	1577	1580	rVB	2570	1760	4.24%	0.310%
13	11.863	1660	1662	1666	rBB	4007	3006	7.25%	0.529%
14	12.557	1777	1780	1783	rBB	11275	9362	22.58%	1.649%
15	12.780	1814	1818	1821	rBB	10558	9812	23.67%	1.728%
16	12.921	1839	1842	1845	rBB	55310	41442	99.95%	7.299%
17	13.110	1871	1874	1877	rVB	1787	1896	4.57%	0.334%
18	13.180	1884	1886	1889	rBB	1382	1003	2.42%	0.177%
19	13.416	1924	1926	1930	rVB	735	974	2.35%	0.172%
20	13.492	1934	1939	1941	rVV	707	1089	2.63%	0.192%
21	13.598	1955	1957	1959	rBV	983	997	2.40%	0.176%
22	13.616	1959	1960	1963	rVB	1885	1424	3.43%	0.251%
23	13.821	1990	1995	1999	rVB3	3960	5561	13.41%	0.979%
24	13.868	1999	2003	2005	rBV2	981	1197	2.89%	0.211%
25	13.933	2009	2014	2015	rVB2	1788	1638	3.95%	0.288%
26	13.980	2018	2022	2025	rBV2	33367	30679	73.99%	5.403%
27	14.010	2025	2027	2030	rVV	3989	3238	7.81%	0.570%
28	14.068	2034	2037	2038	rBV2	1801	1263	3.05%	0.222%
29	14.139	2046	2049	2050	rBV	1294	1317	3.18%	0.232%
30	14.198	2056	2059	2061	rBV2	1252	1109	2.67%	0.195%
31	14.239	2063	2066	2069	rBV2	1165	1653	3.99%	0.291%
32	14.263	2069	2070	2074	rVB2	1134	1272	3.07%	0.224%
33	14.310	2076	2078	2079	rBV	1475	1118	2.70%	0.197%
34	14.439	2097	2100	2103	rBV3	2975	4591	11.07%	0.809%
35	14.468	2103	2105	2107	rVV2	2279	2403	5.80%	0.423%
36	14.492	2107	2109	2111	rVV2	2506	2341	5.65%	0.412%

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37	14.521	2111	2114	2115	rVV2	1699	2003	4.83%	0.353%
38	14.545	2117	2118	2119	rVV	2721	1331	3.21%	0.234%
39	14.580	2119	2124	2125	rVV3	1930	3038	7.33%	0.535%
40	14.621	2127	2131	2132	rVV3	1727	2414	5.82%	0.425%
41	14.674	2136	2140	2142	rVV4	1834	2303	5.55%	0.406%
42	14.704	2144	2145	2148	rVV2	1904	1311	3.16%	0.231%
43	14.757	2150	2154	2156	rVV2	1567	2125	5.13%	0.374%
44	14.792	2158	2160	2162	rVB2	1242	1118	2.70%	0.197%
45	14.815	2162	2164	2167	rBV3	1158	1428	3.44%	0.251%
46	14.845	2167	2169	2171	rVV3	968	1124	2.71%	0.198%
47	14.998	2189	2195	2196	rBV5	2390	3785	9.13%	0.667%
48	15.015	2196	2198	2201	rVB2	4731	4896	11.81%	0.862%
49	15.210	2227	2231	2232	rVV3	1932	1730	4.17%	0.305%
50	15.286	2242	2244	2246	rVV2	1978	1765	4.26%	0.311%
51	15.315	2246	2249	2251	rVV2	3548	3272	7.89%	0.576%
52	15.351	2253	2255	2256	rBV	1855	1105	2.67%	0.195%
53	15.374	2256	2259	2263	rVV2	6071	6490	15.65%	1.143%
54	15.439	2266	2270	2276	rVB2	16509	20628	49.75%	3.633%
55	15.645	2302	2305	2306	rBV3	2121	2422	5.84%	0.427%
56	15.657	2306	2307	2309	rVV2	2296	1920	4.63%	0.338%
57	15.686	2309	2312	2314	rVV3	1676	2177	5.25%	0.383%
58	15.745	2318	2322	2329	rVV5	1597	4204	10.14%	0.740%
59	15.874	2341	2344	2346	rBV3	1173	1178	2.84%	0.207%
60	15.927	2351	2353	2355	rBV2	1851	1598	3.85%	0.281%
61	16.004	2364	2366	2368	rBV	1093	1102	2.66%	0.194%
62	16.098	2379	2382	2384	rBV2	1712	2364	5.70%	0.416%
63	16.115	2384	2385	2388	rVV2	2590	2235	5.39%	0.394%
64	16.145	2388	2390	2395	rVV5	1360	1918	4.63%	0.338%
65	16.210	2400	2401	2404	rVV3	1848	1425	3.44%	0.251%
66	16.268	2407	2411	2413	rBV3	1493	1921	4.63%	0.338%
67	16.304	2415	2417	2420	rVV2	1280	1664	4.01%	0.293%
68	16.333	2420	2422	2425	rVV4	1187	1458	3.52%	0.257%
69	16.374	2425	2429	2433	rVV4	1500	2728	6.58%	0.480%
70	16.415	2433	2436	2439	rVB4	1564	1889	4.56%	0.333%
71	16.527	2452	2455	2456	rBV	952	1164	2.81%	0.205%
72	16.545	2456	2458	2461	rVV3	2015	2063	4.98%	0.363%
73	16.745	2489	2492	2493	rVB	1327	1084	2.61%	0.191%
74	16.768	2493	2496	2498	rBV2	1253	1273	3.07%	0.224%
75	16.792	2498	2500	2503	rVV2	1496	1591	3.84%	0.280%
76	16.821	2503	2505	2506	rVV	1557	1078	2.60%	0.190%

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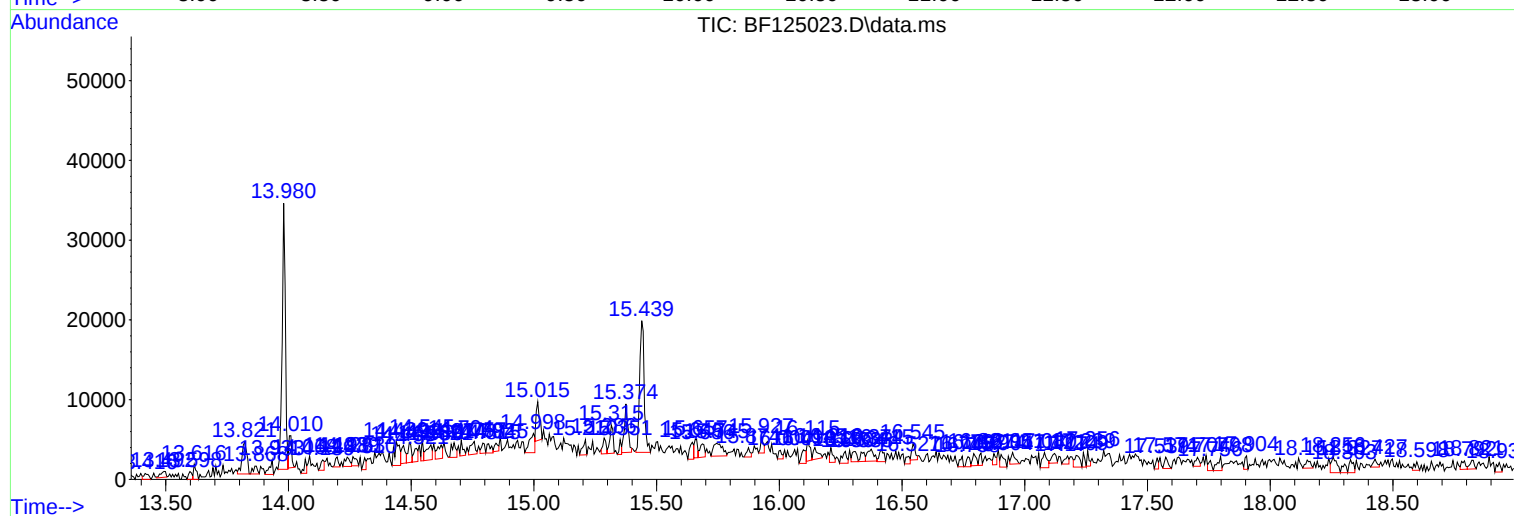
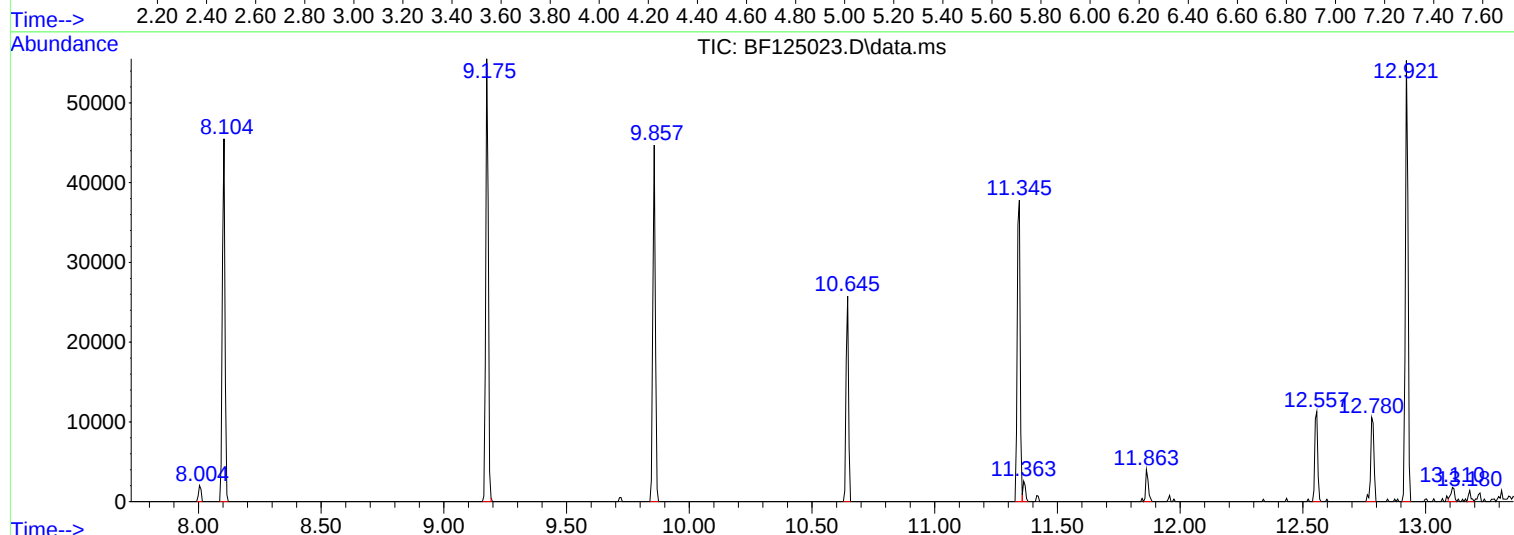
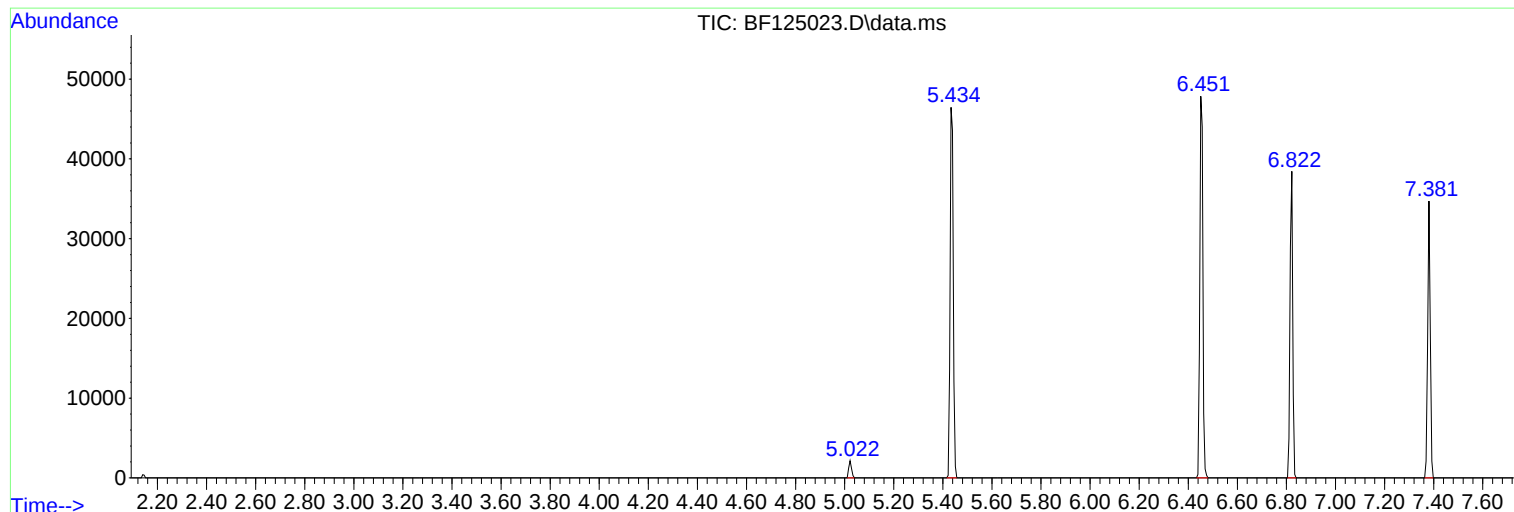
Peak separation: 5

Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	16.874	2513	2514	2516	rVV2	1404	1154	2.78%	0.203%
78	16.904	2518	2519	2523	rVB3	1558	1673	4.04%	0.295%
79	16.951	2525	2527	2532	rVB3	1413	1583	3.82%	0.279%
80	17.080	2546	2549	2552	rBV4	1899	2601	6.27%	0.458%
81	17.121	2552	2556	2558	rVV4	1196	1548	3.73%	0.273%
82	17.174	2561	2565	2566	rBV	1039	1179	2.84%	0.208%
83	17.209	2569	2571	2575	rVB3	1461	2054	4.95%	0.362%
84	17.239	2575	2576	2578	rBV2	1797	1031	2.49%	0.182%
85	17.256	2578	2579	2581	rBV2	1927	1124	2.71%	0.198%
86	17.539	2625	2627	2628	rBV	1354	1012	2.44%	0.178%
87	17.574	2631	2633	2637	rVB3	1434	1403	3.38%	0.247%
88	17.704	2652	2655	2657	rBV3	1138	1288	3.11%	0.227%
89	17.756	2662	2664	2666	rVB2	1258	1010	2.44%	0.178%
90	17.798	2666	2671	2672	rBV3	1752	1934	4.66%	0.341%
91	17.904	2687	2689	2690	rVB2	1685	1104	2.66%	0.194%
92	18.151	2728	2731	2735	rVV3	1014	1411	3.40%	0.249%
93	18.256	2747	2749	2754	rVB3	1973	2844	6.86%	0.501%
94	18.303	2754	2757	2759	rVB2	1084	1228	2.96%	0.216%
95	18.327	2759	2761	2764	rBV2	1638	1989	4.80%	0.350%
96	18.427	2775	2778	2781	rBV3	982	1110	2.68%	0.195%
97	18.598	2804	2807	2809	rVB2	1058	975	2.35%	0.172%
98	18.792	2837	2840	2842	rVV3	1061	1254	3.02%	0.221%
99	18.821	2842	2845	2848	rVB2	1107	1269	3.06%	0.223%
100	18.933	2861	2864	2866	rBV2	1185	1277	3.08%	0.225%

Sum of corrected areas: 567806

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P



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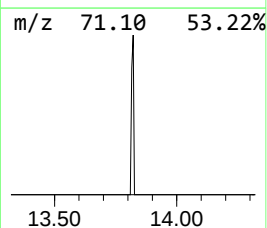
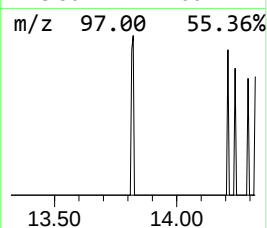
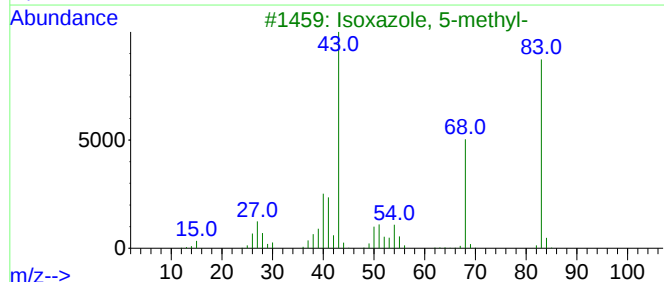
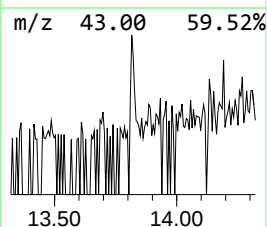
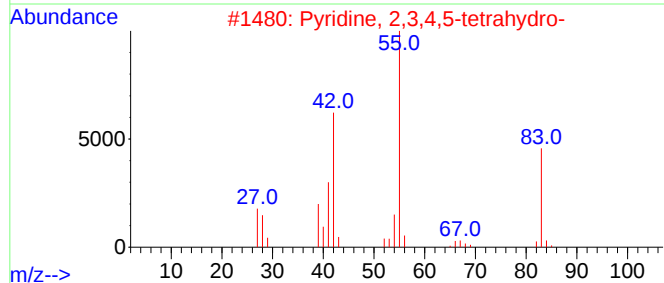
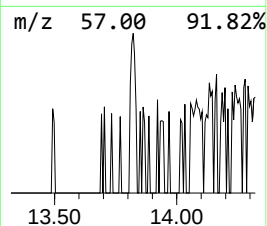
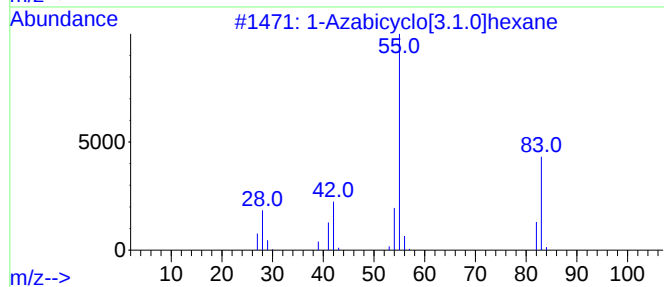
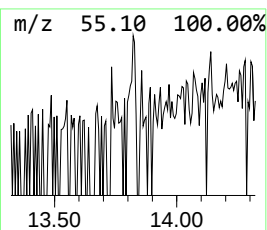
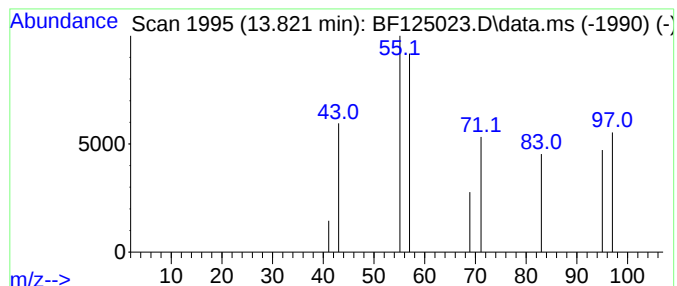
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 1 unknown13.821 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.821	3.63 ng	5561	Chrysene-d12	13.980

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Azabicyclo[3.1.0]hexane	83	C5H9N	000285-76-7	5
2			Pyridine, 2,3,4,5-tetrahydro-	83	C5H9N	000505-18-0	4
3			Isoxazole, 5-methyl-	83	C4H5NO	005765-44-6	4
4			1H-Pyrrole-2,5-dione	97	C4H3NO2	000541-59-3	4
5			Azetidine, 2-methyl-	71	C4H9N	019812-49-8	3



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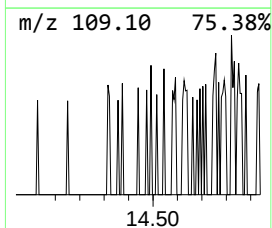
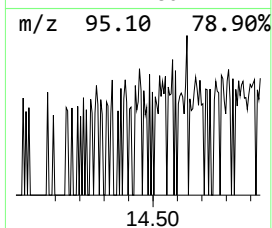
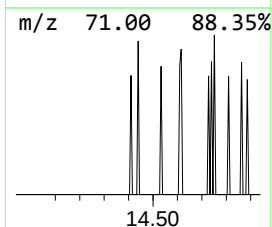
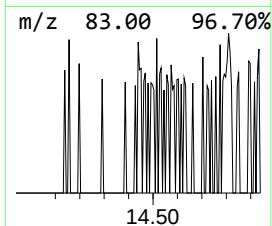
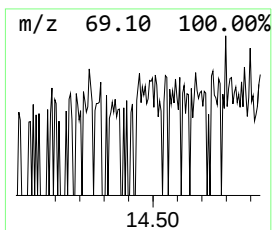
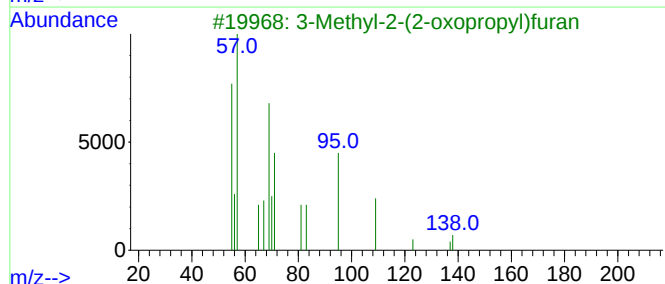
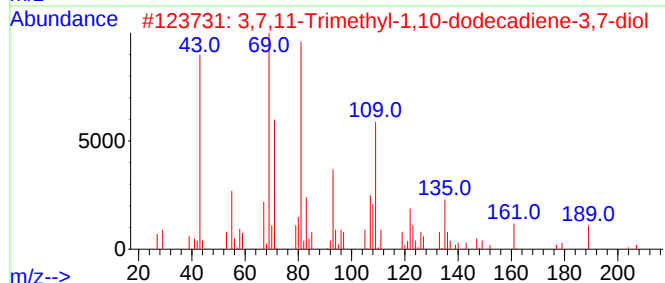
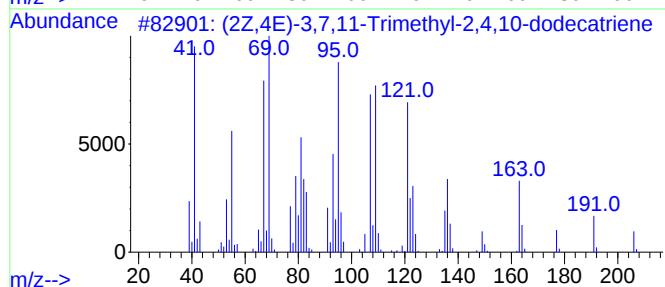
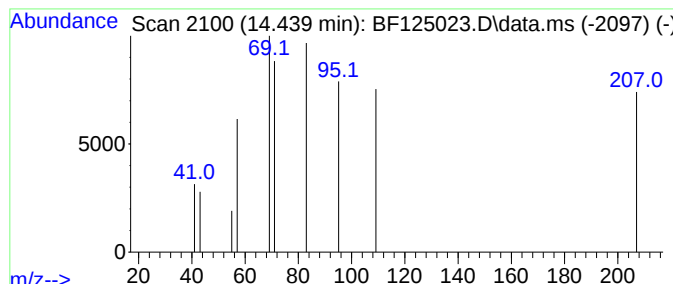
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 unknown14.439 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.439	2.99 ng	4591	Chrysene-d12	13.980

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	(2Z,4E)-3,7,11-Trimethyl-2,4,10-...	206	C15H26	172549-29-0	9		
2	3,7,11-Trimethyl-1,10-dodecadien...	240	C15H28O2	1000432-46-6	4		
3	3-Methyl-2-(2-oxopropyl)furan	138	C8H10O2	087773-62-4	4		
4	1-Tridecyn-4-ol	196	C13H24O	074646-37-0	4		
5	4-t-Butyl-2-(1-methyl-2-nitroeth...	241	C13H23NO3	1000192-74-8	4		



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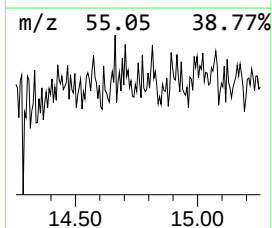
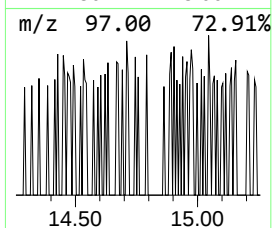
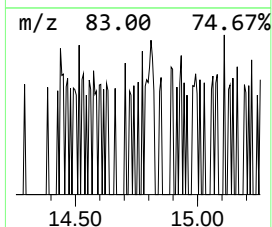
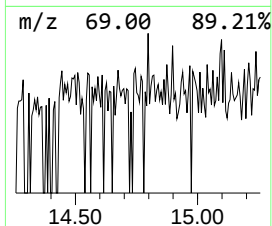
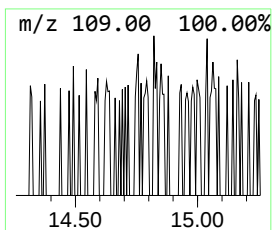
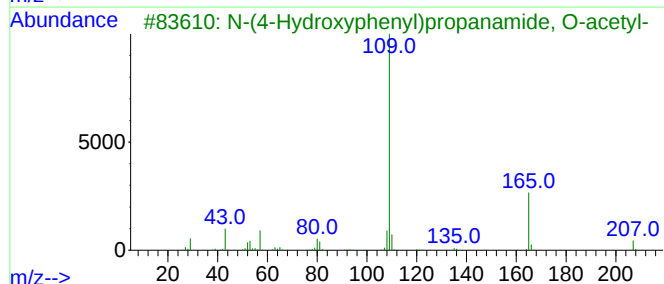
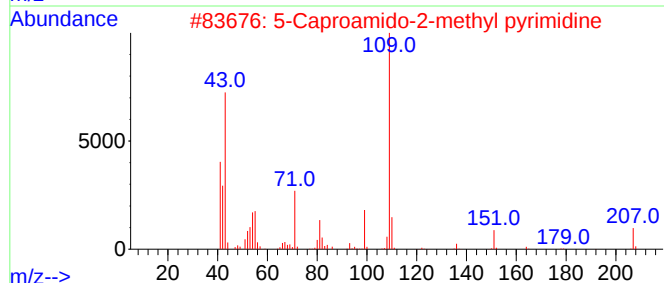
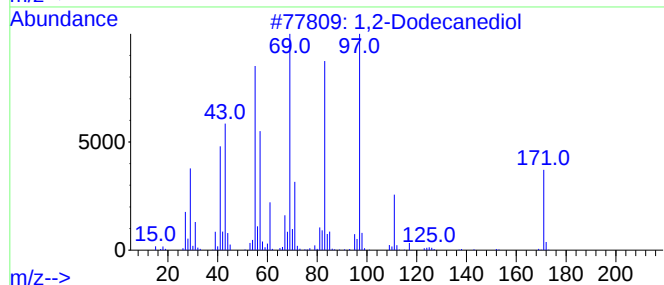
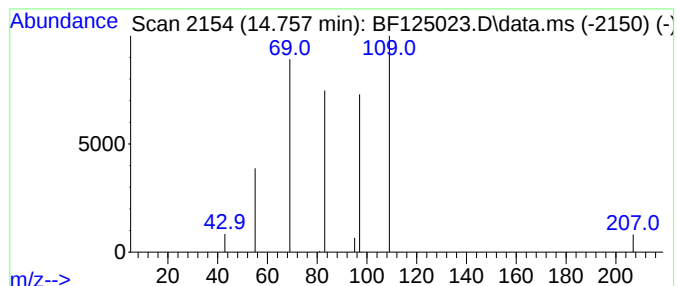
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown14.757 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.757	2.06 ng	2125	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Dodecanediol	202	C12H26O2	001119-87-5	36
2			5-Caproamido-2-methyl pyrimidine	207	C11H17N3O	1000213-95-8	4
3			N-(4-Hydroxyphenyl)propanamide, ...	207	C11H13NO3	1824295-97-7	4
4			Cyclopropane carboxamide, 2-cycl...	207	C13H21NO	331416-19-4	4
5			1H-Imidazol-2-amine	83	C3H5N3	007720-39-0	4



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Data File : BF125023.D
Acq On : 10 Aug 2021 2:59
Operator : JU/CG
Sample : M3314-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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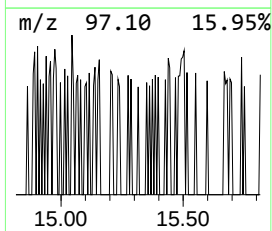
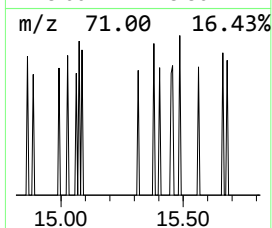
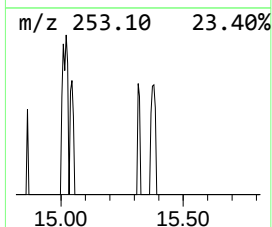
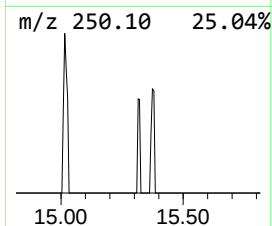
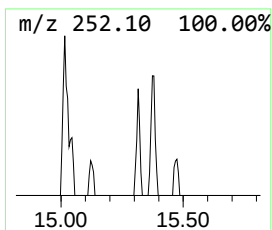
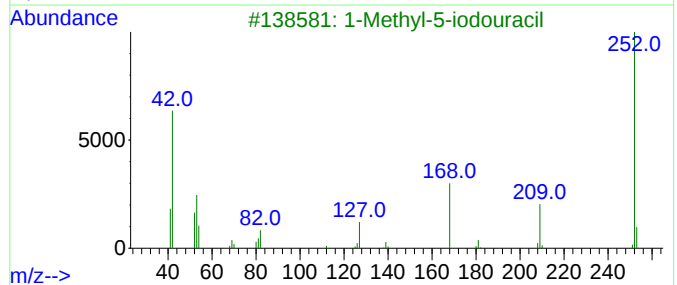
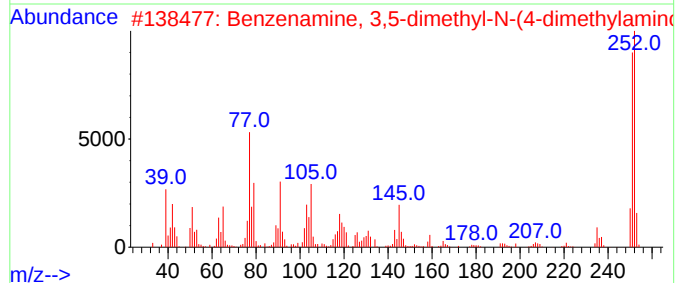
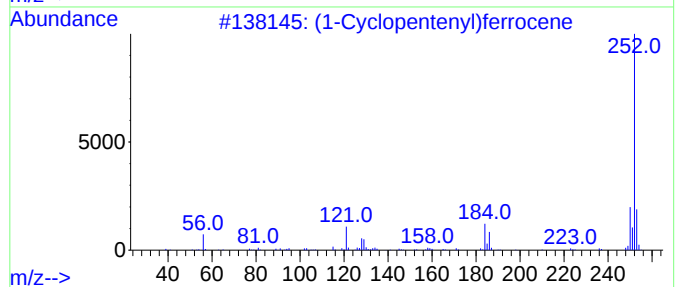
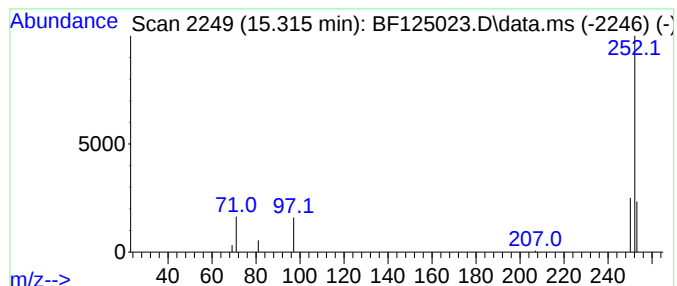
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown15.316 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.316	3.17 ng	3272	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	9
2			Benzenamine, 3,5-dimethyl-N-(4-d...	252	C17H20N2	303769-91-7	9
3			1-Methyl-5-iodouracil	252	C5H5IN2O2	045774-47-8	5
4			Benzo[j]fluoranthene	252	C20H12	000205-82-3	5
5			Benzo[k]fluoranthene	252	C20H12	000207-08-9	5



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
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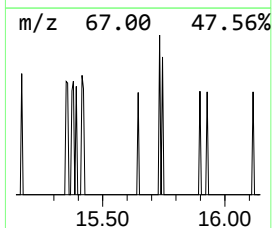
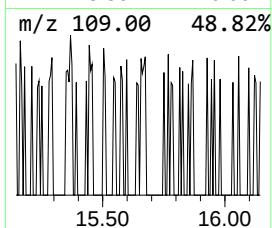
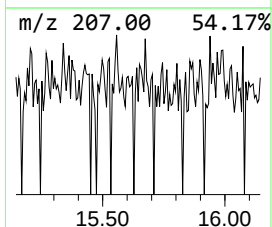
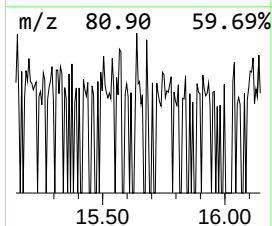
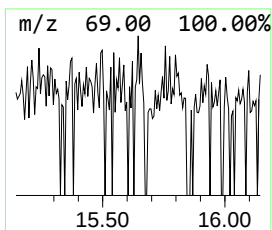
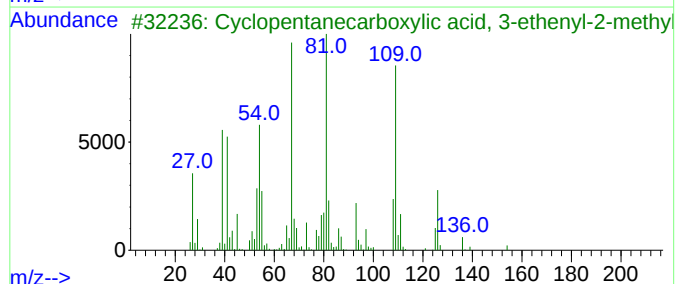
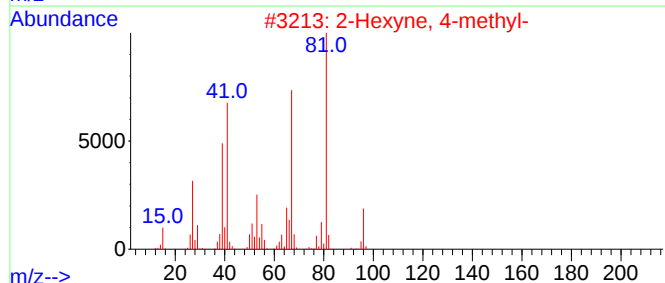
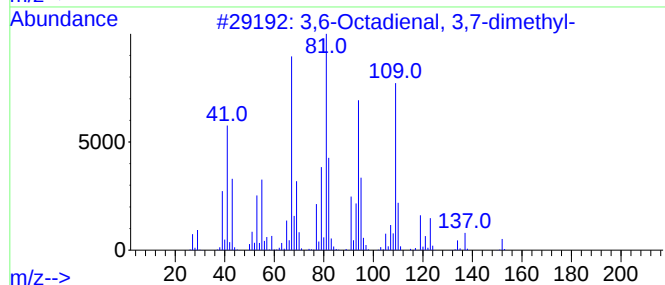
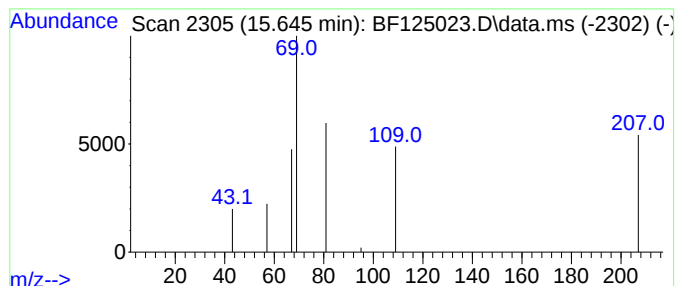
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 unknown15.645 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.645	2.35 ng	2422	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,6-Octadienal, 3,7-dimethyl-	152	C10H16O	055722-59-3	9
2			2-Hexyne, 4-methyl-	96	C7H12	020198-49-6	4
3			Cyclopentanecarboxylic acid, 3-e...	154	C9H14O2	108451-44-1	4
4			Cyclohexanone, 2-(1-mercapto-1-m...	186	C10H18OS	033281-91-3	4
5			Cyclopropane, 1,1-dimethyl-2-(2-...	124	C9H16	069147-03-1	4



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
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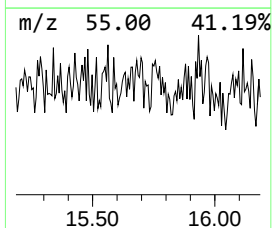
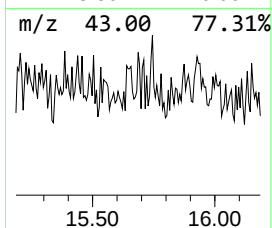
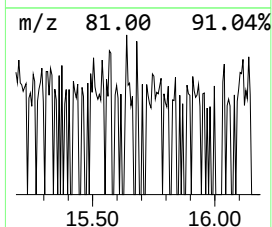
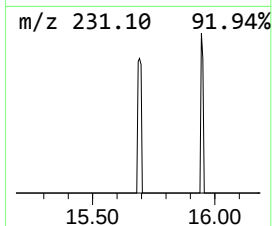
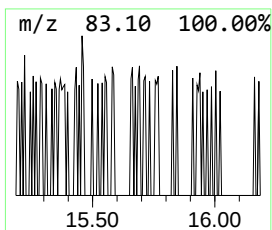
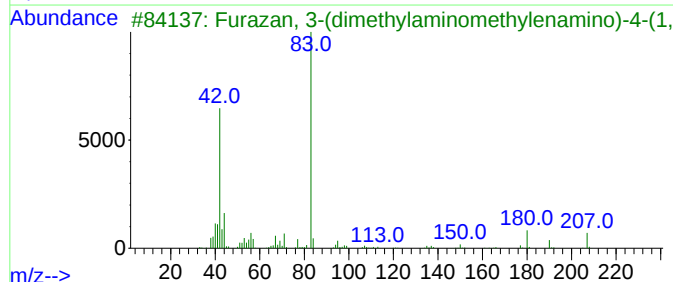
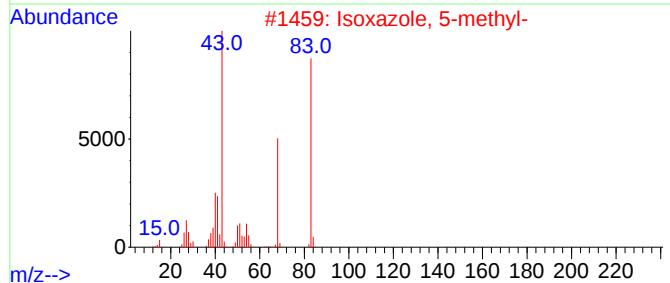
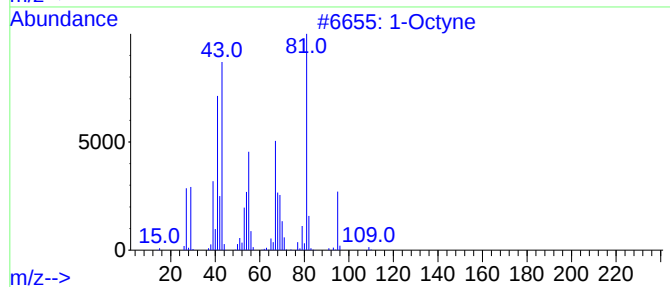
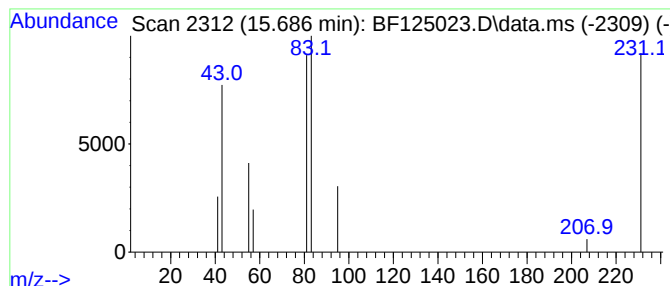
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 unknown15.686 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.686	2.11 ng	2177	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Octyne	110	C8H14	000629-05-0	9
2			Isoxazole, 5-methyl-	83	C4H5NO	005765-44-6	4
3			Furazan, 3-(dimethylaminomethyle...	207	C7H9N7O	294667-75-7	4
4			N-(2-Methoxyphenyl)-benzene-sulp...	231	C13H13NOS	111685-76-8	2
5			Hexanal trans-2-hexenyl hexyl ac...	284	C18H36O2	1000430-94-8	2



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
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Misc :
ALS Vial : 20 Sample Multiplier: 1

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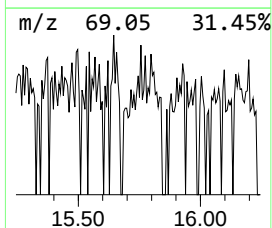
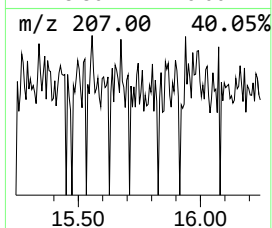
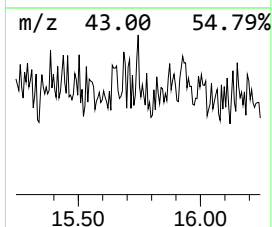
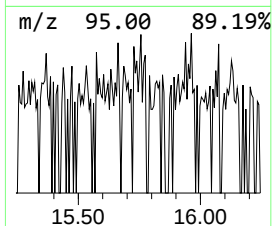
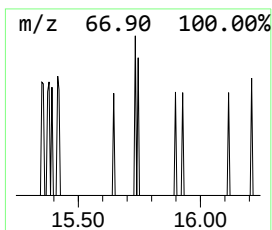
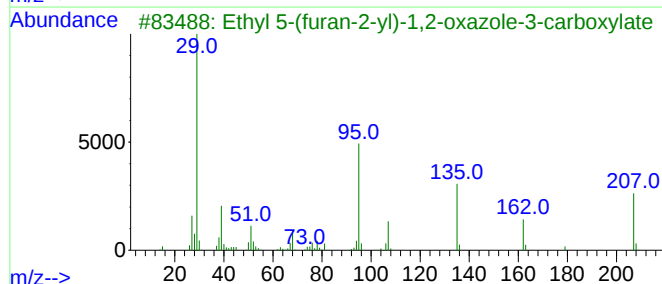
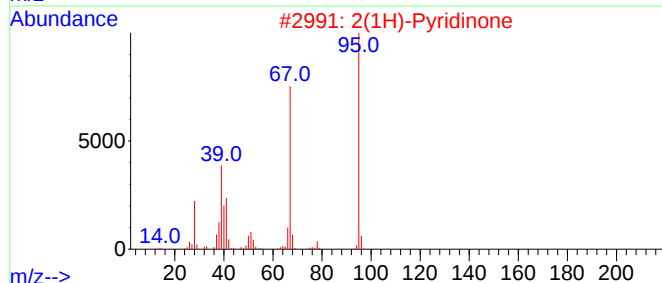
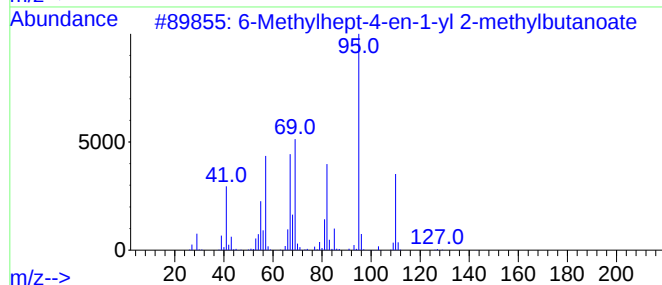
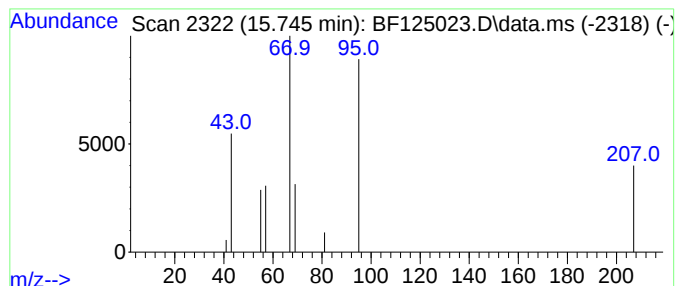
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 unknown15.745 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.745	4.08 ng	4204	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methylhept-4-en-1-yl 2-methylb...	212	C13H24O2	1215128-05-4	9		
2	2(1H)-Pyridinone	95	C5H5NO	000142-08-5	5		
3	Ethyl 5-(furan-2-yl)-1,2-oxazole...	207	C10H9NO4	1000411-40-9	4		
4	Chlorine dioxide	67	ClO2	010049-04-4	4		
5	4-Methyl-1,3-heptadiene	110	C8H14	017603-57-5	4		



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
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Misc :
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Instrument :
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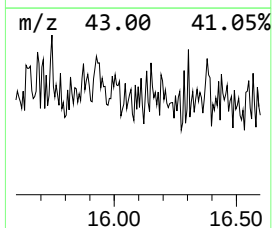
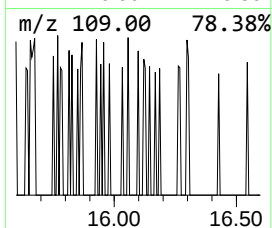
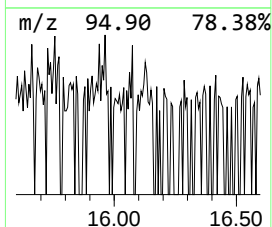
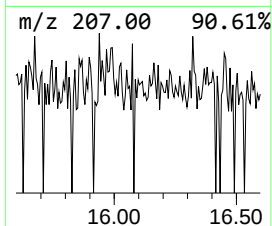
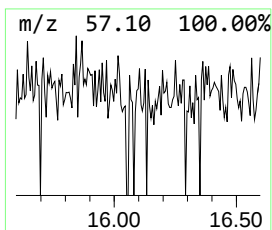
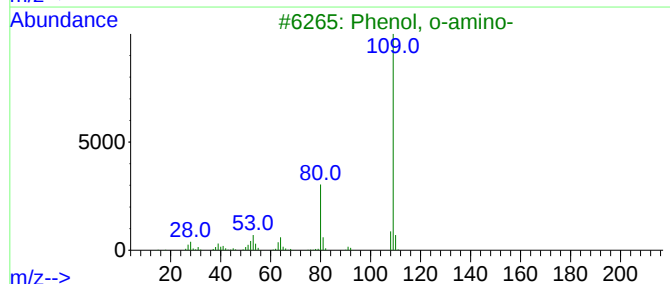
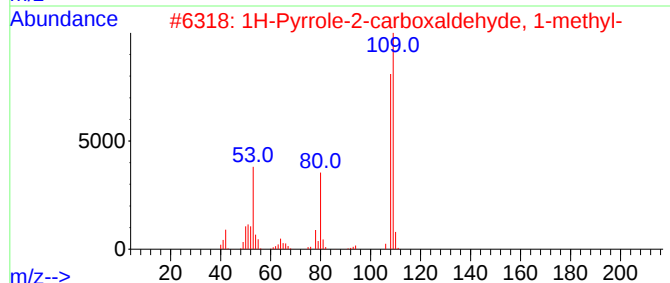
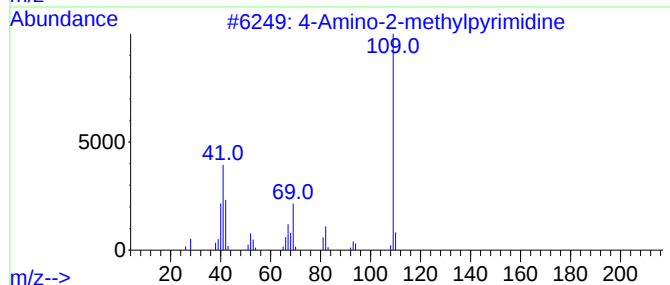
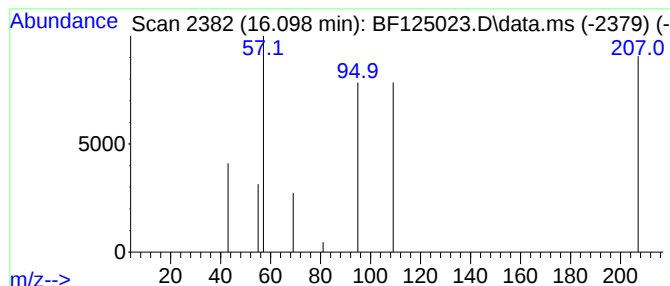
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 4-Amino-2-methylpyrimidine Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.098	2.29 ng	2364	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Amino-2-methylpyrimidine	109	C5H7N3	000074-69-1	4
2			1H-Pyrrole-2-carboxaldehyde, 1-m...	109	C6H7NO	001192-58-1	4
3			Phenol, o-amino-	109	C6H7NO	000095-55-6	4
4			1H-Imidazole, 2-ethyl-4-methyl-	110	C6H10N2	000931-36-2	4
5			2-Amino-4-methylpyrimidine	109	C5H7N3	000108-52-1	4



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
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Misc :
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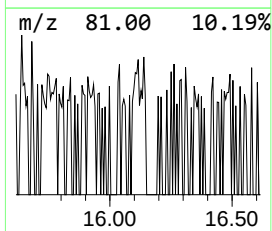
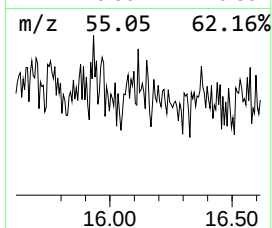
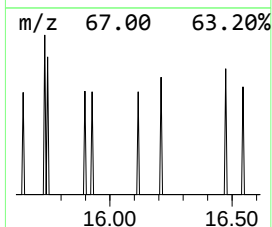
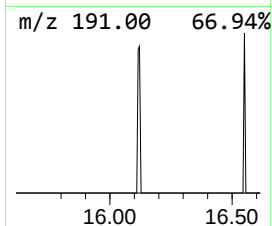
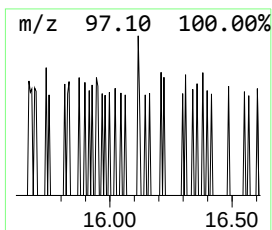
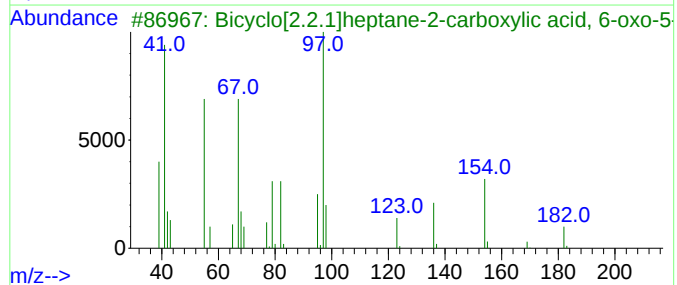
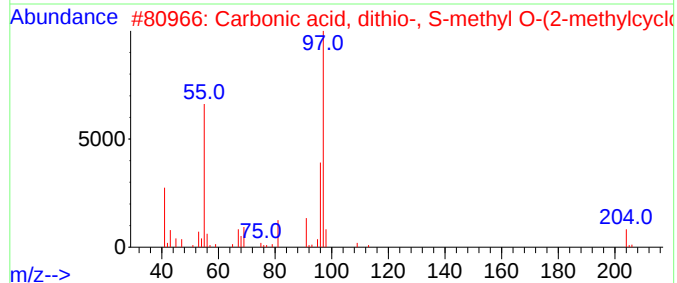
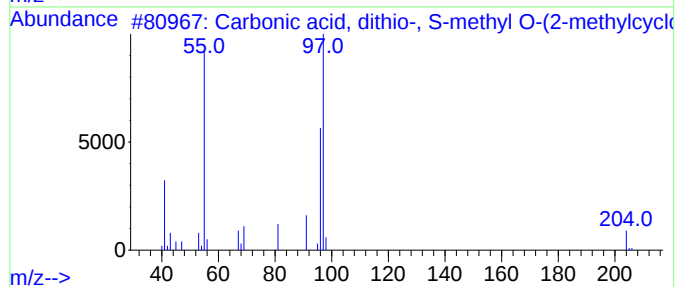
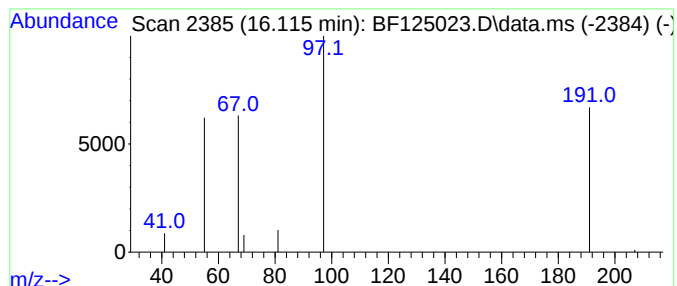
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 unknown16.115 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.115	2.17 ng	2235	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, dithio-, S-methyl...	204	C9H16O5S2	015288-13-8	9
2			Carbonic acid, dithio-, S-methyl...	204	C9H16O5S2	015288-12-7	9
3			Bicyclo[2.2.1]heptane-2-carboxyl...	210	C11H14O4	111509-63-8	9
4			Oxazole, 4,5-dimethyl-	97	C5H7NO	020662-83-3	7
5			4H-1,2,4-Triazole, 4-ethyl-	97	C4H7N3	043183-55-7	5



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
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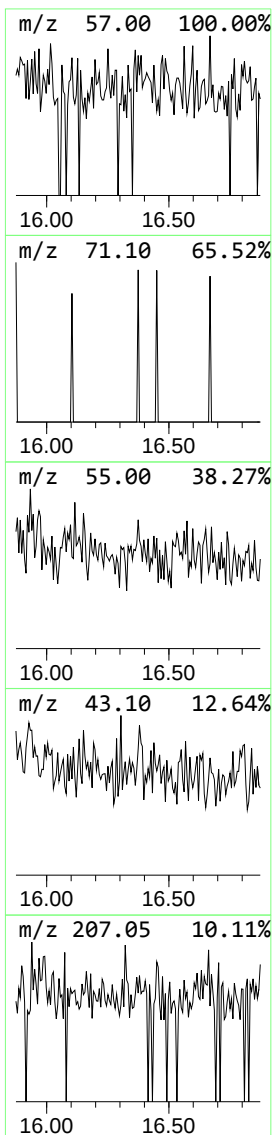
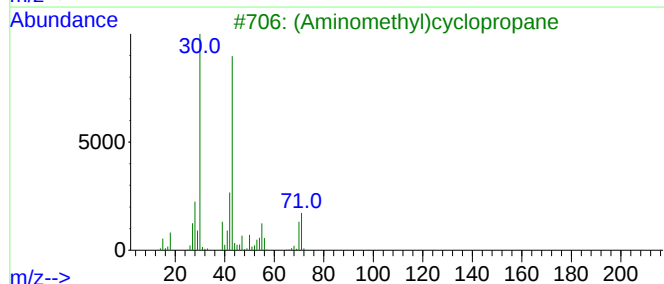
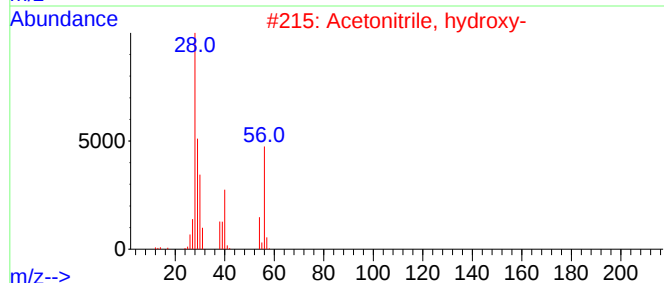
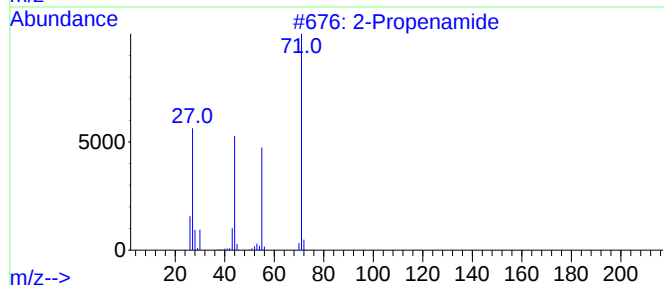
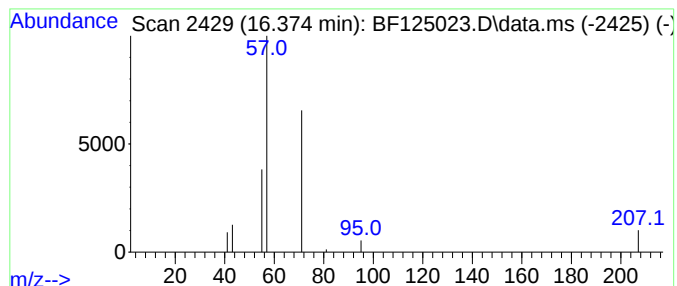
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 unknown16.374 Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.374	2.64 ng	2728	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Propenamide	71	C3H5NO	000079-06-1	5
2			Acetonitrile, hydroxy-	57	C2H3NO	000107-16-4	3
3			(Aminomethyl)cyclopropane	71	C4H9N	002516-47-4	3
4			Azetidine	57	C3H7N	000503-29-7	3
5			Oxirane, 2,2'-[1,4-butanediylbis...	202	C10H18O4	002425-79-8	2



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
Acq On : 10 Aug 2021 2:59
Operator : JU/CG
Sample : M3314-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

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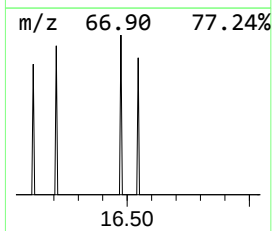
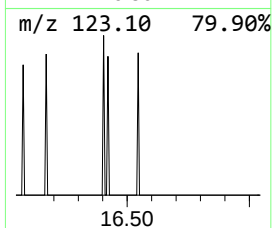
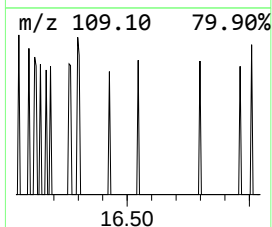
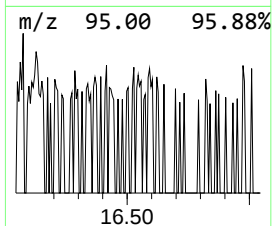
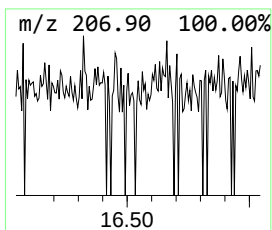
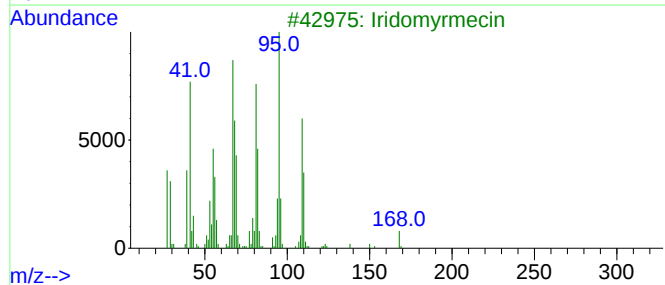
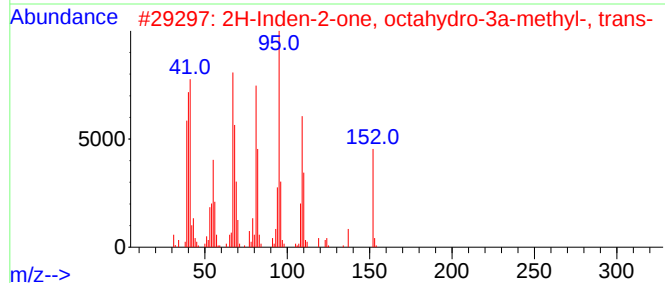
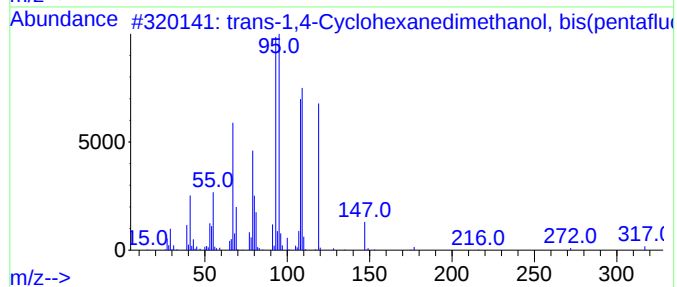
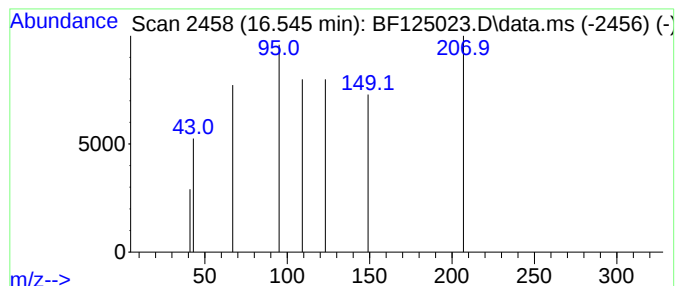
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 unknown16.545 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.545	2.00 ng	2063	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-1,4-Cyclohexanedimethanol,...	436	C14H14F10O4	1010384-27-9	9
2			2H-Inden-2-one, octahydro-3a-met...	152	C10H16O	020379-99-1	9
3			Iridomyrmecin	168	C10H16O2	000485-43-8	9
4			8-Thiabicyclo[4.3.0]nonane S-oxide	158	C8H14OS	1010215-10-0	9
5			Ethyl 2-octynoate	168	C10H16O2	010519-20-7	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
Acq On : 10 Aug 2021 2:59
Operator : JU/CG
Sample : M3314-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
72-11931

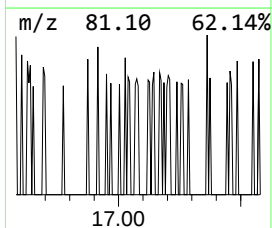
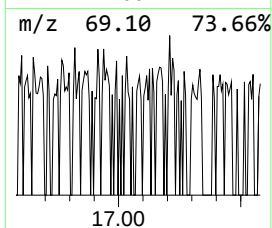
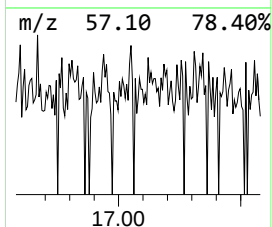
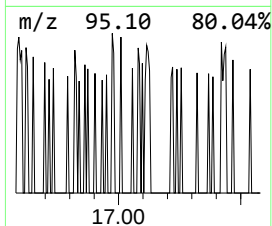
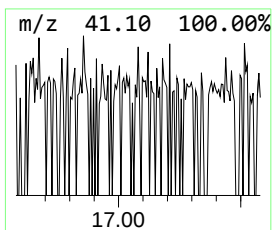
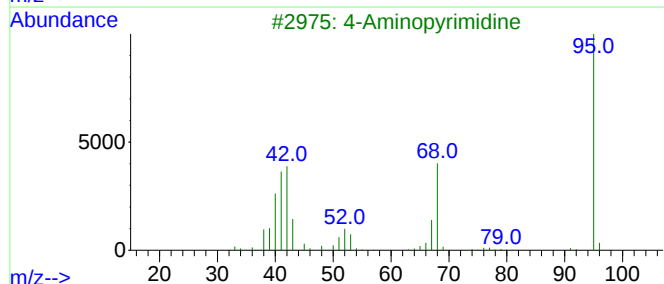
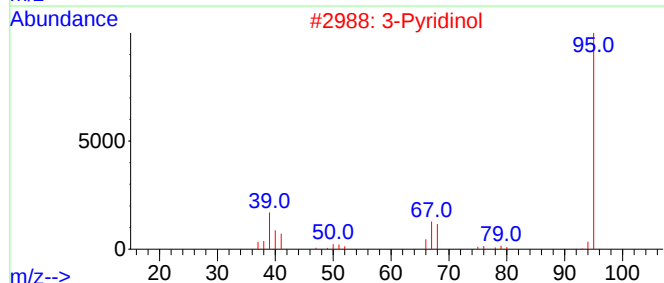
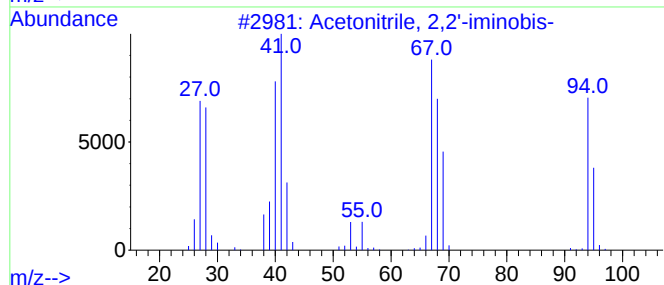
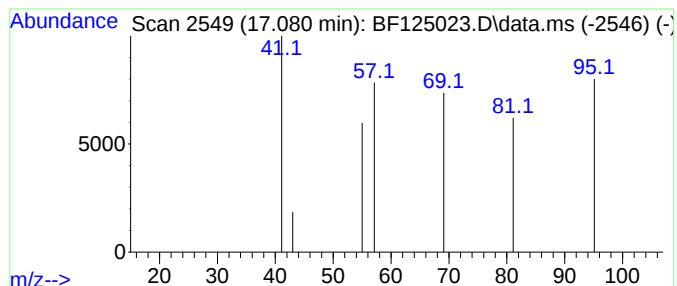
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 unknown17.080 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.080	2.52 ng	2601	Perylene-d12	15.439

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Acetonitrile, 2,2'-iminobis-	95	C4H5N3	000628-87-5	4
2		3-Pyridinol	95	C5H5NO	000109-00-2	2
3		4-Aminopyrimidine	95	C4H5N3	000591-54-8	2
4		Hexanal, 5,5-dimethyl-	128	C8H16O	055320-58-6	2
5		Oxazole	69	C3H3NO	000288-42-6	2



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
Acq On : 10 Aug 2021 2:59
Operator : JU/CG
Sample : M3314-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
72-11931

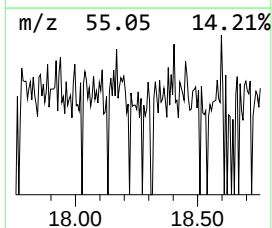
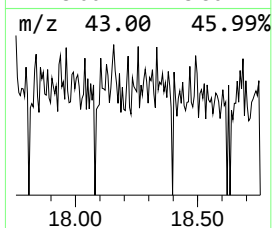
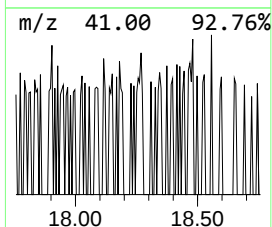
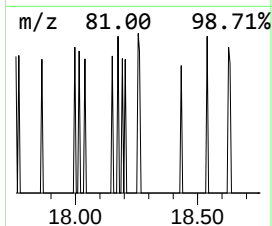
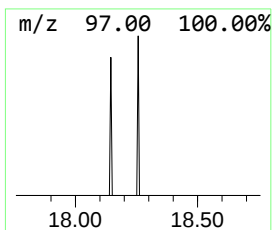
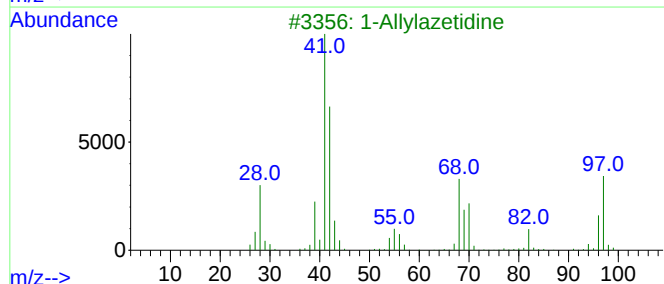
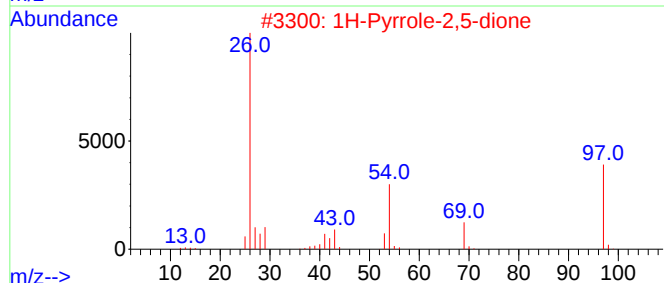
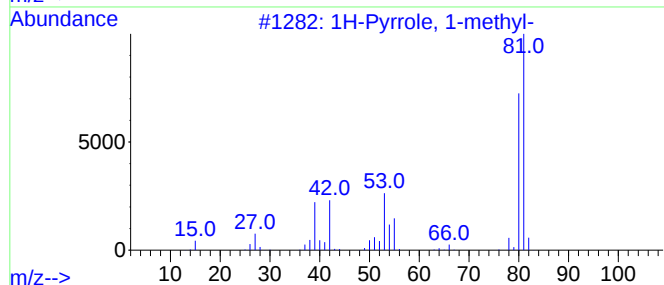
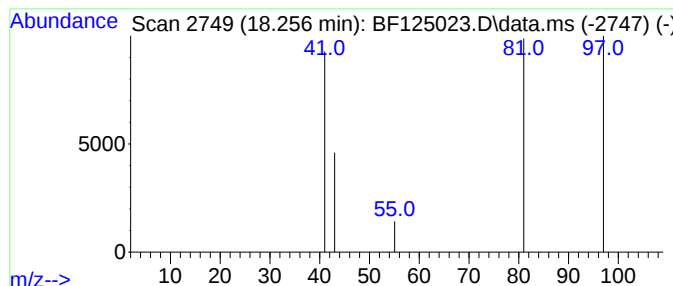
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 unknown18.256 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.256	2.76 ng	2844	Perylene-d12	15.439

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Pyrrole, 1-methyl-	81	C5H7N	000096-54-8	2
2			1H-Pyrrole-2,5-dione	97	C4H3NO2	000541-59-3	2
3			1-Allylazetidine	97	C6H11N	1000195-02-5	2
4			2-Octynoic acid	140	C8H12O2	005663-96-7	1
5			1-Hexyne, 5-methyl-	96	C7H12	002203-80-7	1



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF080921\
Data File : BF125023.D
Acq On : 10 Aug 2021 2:59
Operator : JU/CG
Sample : M3314-01 2X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
72-11931

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF080421.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST0.L
TIC Integration Parameters: LSCINT.P

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unknown14.439	14.439	3.0	ng	4591	5	13.980	30679	20.0
unknown14.757	14.757	2.1	ng	2125	6	15.439	20628	20.0
unknown15.316	15.316	3.2	ng	3272	6	15.439	20628	20.0
unknown15.645	15.645	2.4	ng	2422	6	15.439	20628	20.0
unknown15.686	15.686	2.1	ng	2177	6	15.439	20628	20.0
unknown15.745	15.745	4.1	ng	4204	6	15.439	20628	20.0
4-Amino-2-methy...	16.098	2.3	ng	2364	6	15.439	20628	20.0
unknown16.115	16.115	2.2	ng	2235	6	15.439	20628	20.0
unknown16.374	16.374	2.6	ng	2728	6	15.439	20628	20.0
unknown16.545	16.545	2.0	ng	2063	6	15.439	20628	20.0
unknown17.080	17.080	2.5	ng	2601	6	15.439	20628	20.0
unknown18.256	18.256	2.8	ng	2844	6	15.439	20628	20.0