

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SEEP-1

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_P\Methods\8270E-BP102925.M
 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP026099.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.226	43	49	60	rVB	215739	294513	1.81%	0.155%
2	3.490	86	94	107	rVB2	124824	309144	1.90%	0.163%
3	3.831	144	152	153	rBV2	297510	496117	3.05%	0.261%
4	3.920	165	167	174	rVB2	838279	906028	5.57%	0.477%
5	4.384	237	246	253	rBV2	144897	303548	1.87%	0.160%
6	4.596	278	282	287	rVB	114473	164010	1.01%	0.086%
7	4.737	299	306	317	rVV3	142151	289899	1.78%	0.153%
8	4.831	318	322	329	rVB	110934	175409	1.08%	0.092%
9	5.220	374	388	391	rBV	593420	1540728	9.47%	0.811%
10	5.290	395	400	412	rVB2	3060224	5448344	33.48%	2.867%
11	5.972	508	516	522	rBV	127782	209390	1.29%	0.110%
12	6.102	531	538	548	rVB	138114	288883	1.78%	0.152%
13	6.849	633	665	680	rBV3	3447183	12990120	79.82%	6.835%
14	7.008	685	692	700	rBV	182738	342610	2.11%	0.180%
15	7.625	786	797	800	rBV	141541	237398	1.46%	0.125%
16	7.672	800	805	812	rVB	1123546	1771075	10.88%	0.932%
17	8.284	897	909	921	rBV	555896	1329444	8.17%	0.700%
18	8.431	921	934	937	rBV	909658	1714587	10.54%	0.902%
19	8.461	937	939	944	rVB	418224	516238	3.17%	0.272%
20	8.766	985	991	993	rBV	130040	200465	1.23%	0.105%
21	8.813	993	999	1004	rVV	2959355	4842226	29.75%	2.548%
22	8.866	1004	1008	1016	rVB	367495	591161	3.63%	0.311%
23	9.078	1037	1044	1052	rBV	383861	640639	3.94%	0.337%
24	9.272	1071	1077	1085	rVB	161743	266590	1.64%	0.140%
25	9.378	1085	1095	1098	rBV5	89037	229521	1.41%	0.121%
26	9.455	1104	1108	1116	rVB5	89365	194049	1.19%	0.102%
27	9.708	1144	1151	1159	rVB	205354	450117	2.77%	0.237%
28	9.790	1159	1165	1173	rVB2	239155	362880	2.23%	0.191%
29	9.890	1173	1182	1186	rBV	629568	1396749	8.58%	0.735%
30	9.925	1186	1188	1195	rVV	240652	344154	2.11%	0.181%
31	9.996	1195	1200	1206	rVB3	100071	184943	1.14%	0.097%
32	10.184	1228	1232	1236	rVB	205559	274650	1.69%	0.145%
33	10.355	1256	1261	1269	rBV2	114324	219829	1.35%	0.116%
34	10.443	1269	1276	1284	rVB	1503328	2584240	15.88%	1.360%
35	10.719	1311	1323	1325	rBV7	77515	207139	1.27%	0.109%
36	10.990	1362	1369	1377	rBV	546440	1149935	7.07%	0.605%

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37	12.102	1549	1558	1565	rBV	1124704	1841731	11.32%	0.969%
38	12.278	1579	1588	1594	rBV	1133546	2206939	13.56%	1.161%
39	12.672	1651	1655	1662	rVB4	108887	187890	1.15%	0.099%
40	12.825	1675	1681	1691	rBV10	111504	274854	1.69%	0.145%
41	12.925	1691	1698	1706	rBV	7203255	10746210	66.03%	5.655%
42	13.096	1723	1727	1732	rBV	160467	279163	1.72%	0.147%
43	13.219	1743	1748	1753	rBV	295665	446664	2.74%	0.235%
44	13.272	1753	1757	1761	rBV2	197295	317732	1.95%	0.167%
45	13.537	1797	1802	1810	rVB5	88417	217529	1.34%	0.114%
46	13.978	1873	1877	1881	rBV4	137581	208326	1.28%	0.110%
47	14.113	1896	1900	1904	rBV2	164525	247252	1.52%	0.130%
48	14.237	1917	1921	1926	rBV3	379684	591992	3.64%	0.312%
49	14.313	1926	1934	1947	rVB	2388982	4165029	25.59%	2.192%
50	14.990	2042	2049	2054	rBV4	168484	352288	2.16%	0.185%
51	15.084	2055	2065	2070	rBV2	1795598	3272615	20.11%	1.722%
52	15.401	2111	2119	2131	rBV	4222388	6999864	43.01%	3.683%
53	15.701	2165	2170	2180	rBV	1245846	1894811	11.64%	0.997%
54	15.831	2186	2192	2202	rVB2	6778875	9886331	60.75%	5.202%
55	16.019	2216	2224	2230	rVB2	360656	772787	4.75%	0.407%
56	16.278	2265	2268	2274	rVB	301579	440288	2.71%	0.232%
57	16.372	2280	2284	2288	rBV	387474	557228	3.42%	0.293%
58	16.537	2308	2312	2317	rBV2	461234	684117	4.20%	0.360%
59	16.631	2322	2328	2341	rVB	1847378	3255606	20.01%	1.713%
60	16.760	2347	2350	2355	rVB	228577	346020	2.13%	0.182%
61	16.931	2372	2379	2385	rVB	1246306	1892413	11.63%	0.996%
62	17.131	2407	2413	2419	rVB2	2854666	4404487	27.06%	2.318%
63	17.213	2420	2427	2435	rVV3	962463	1686178	10.36%	0.887%
64	17.284	2435	2439	2444	rVB	265430	409648	2.52%	0.216%
65	17.378	2452	2455	2464	rBV3	410246	748498	4.60%	0.394%
66	17.548	2480	2484	2489	rBV3	363095	533994	3.28%	0.281%
67	17.648	2496	2501	2512	rVB	2081437	3044802	18.71%	1.602%
68	17.801	2524	2527	2530	rBV4	113831	186491	1.15%	0.098%
69	18.095	2572	2577	2584	rBV	2247876	2970888	18.26%	1.563%
70	18.301	2608	2612	2621	rBV	2032511	3093552	19.01%	1.628%
71	19.160	2755	2758	2764	rBV2	194866	295226	1.81%	0.155%
72	19.360	2790	2792	2798	rVB	291179	316915	1.95%	0.167%
73	19.419	2799	2802	2807	rBV	606849	820028	5.04%	0.432%
74	19.654	2838	2842	2858	rBV2	1170149	2331799	14.33%	1.227%
75	19.872	2873	2879	2885	rBV	12152536	16273825	100.00%	8.563%
76	20.054	2908	2910	2914	rBV	215140	279019	1.71%	0.147%

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 Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

77	20.825	3037	3041	3045	rVB	1812063	2077370	12.77%	1.093%
78	20.878	3047	3050	3054	rBV2	271738	550875	3.39%	0.290%
79	21.272	3113	3117	3125	rVB	804692	1251286	7.69%	0.658%
80	21.577	3164	3169	3178	rVB	3485646	5252180	32.27%	2.764%
81	23.848	3549	3555	3573	rVB2	817573	2161184	13.28%	1.137%
82	24.818	3710	3720	3728	rBV	3510925	8519046	52.35%	4.483%
83	24.913	3728	3736	3750	rVB	2178561	5806114	35.68%	3.055%
84	25.971	3906	3916	3931	rVB4	3414422	10854473	66.70%	5.712%
85	27.348	4137	4150	4167	rVB2	3842807	13916316	85.51%	7.323%
86	28.995	4417	4430	4451	rVB2	1443074	6704161	41.20%	3.528%

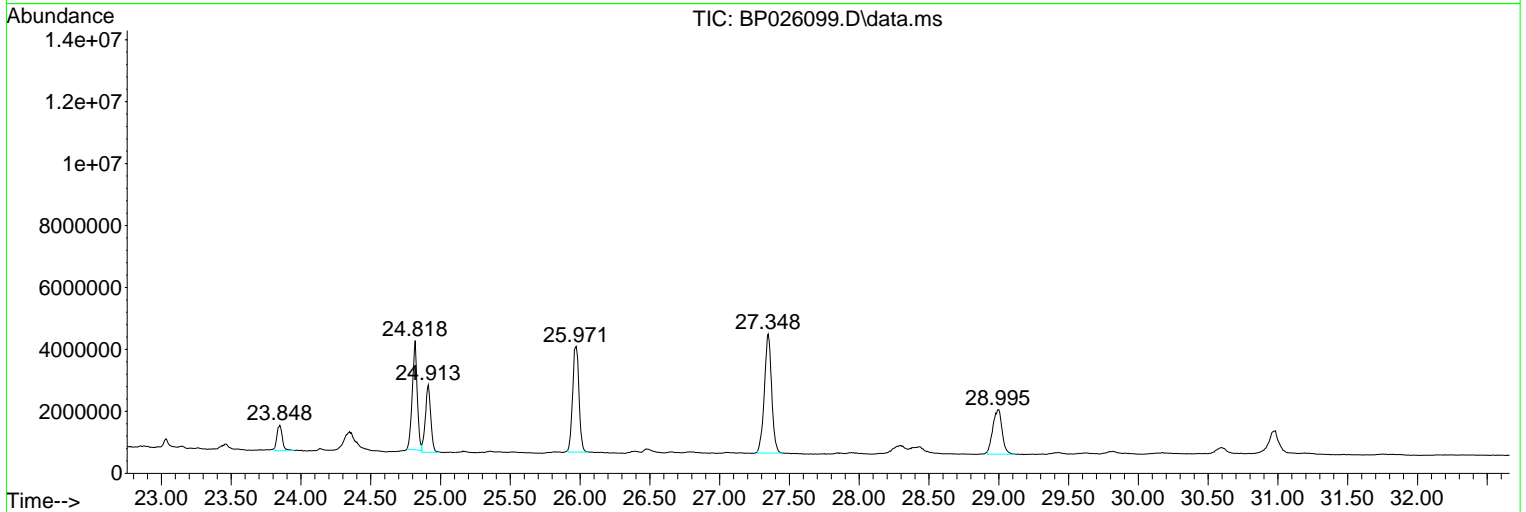
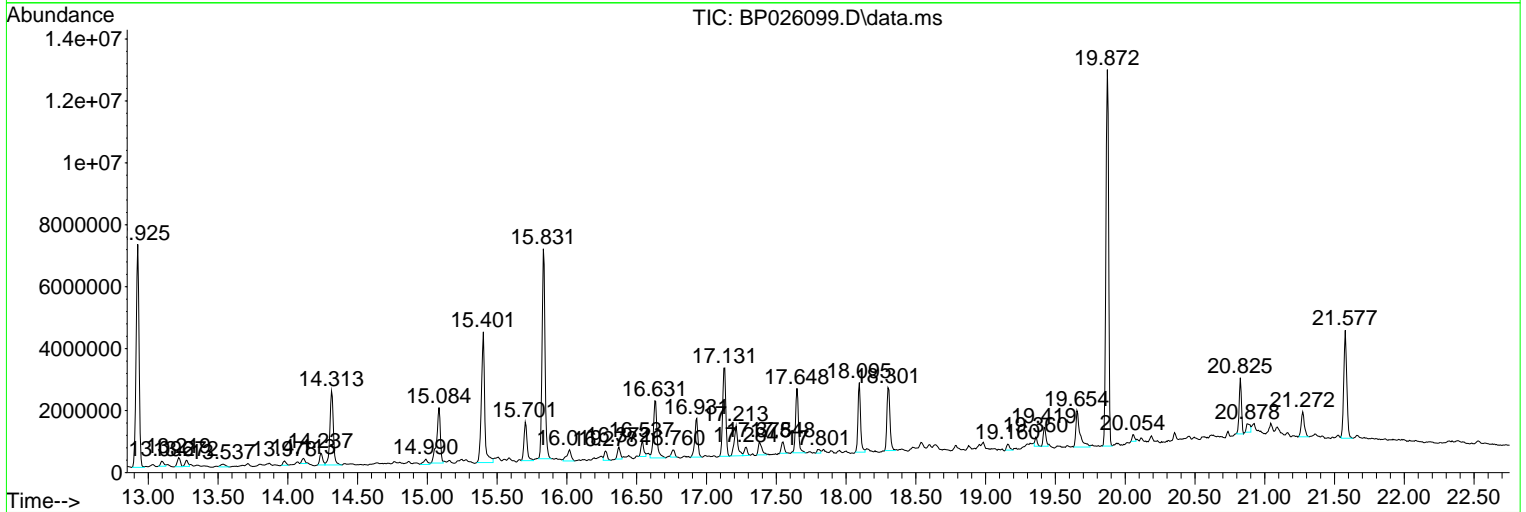
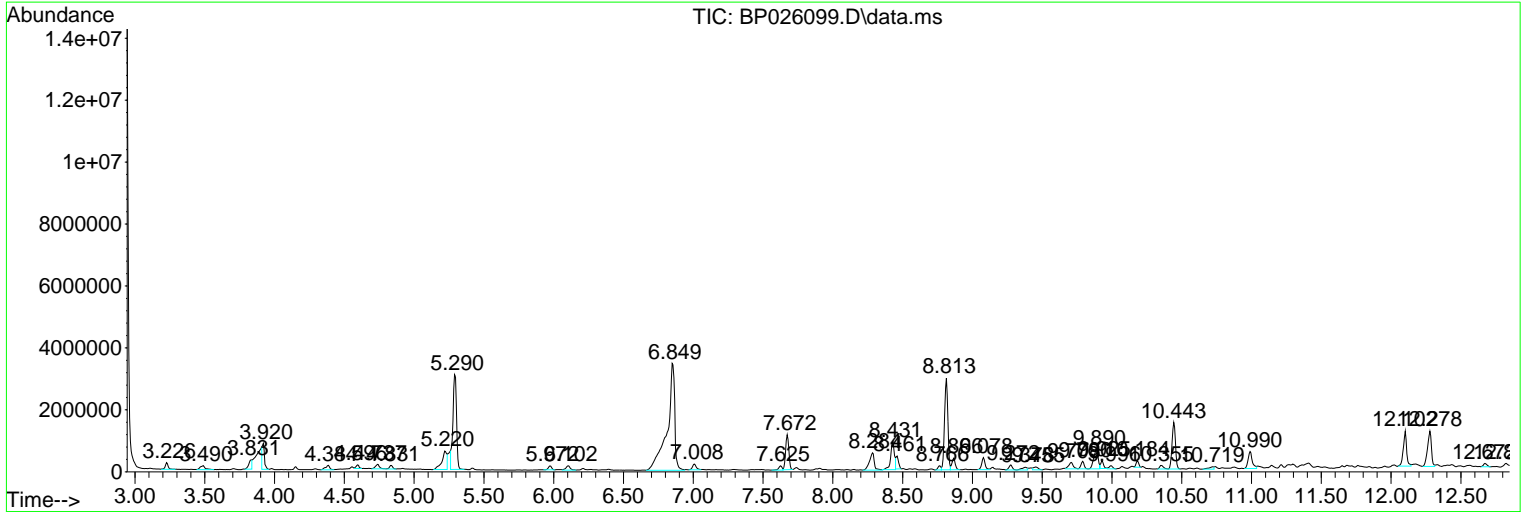
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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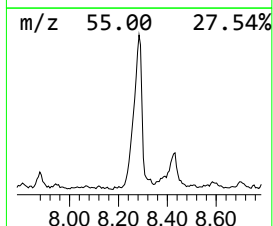
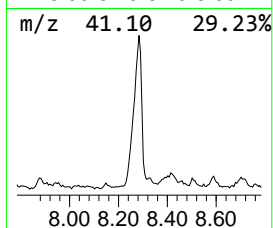
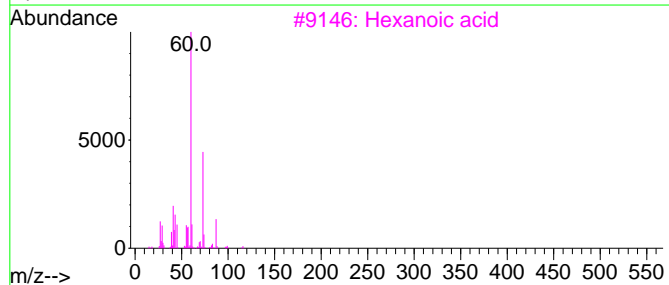
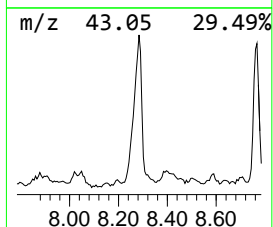
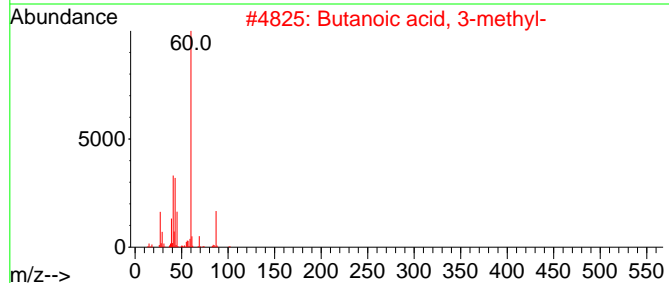
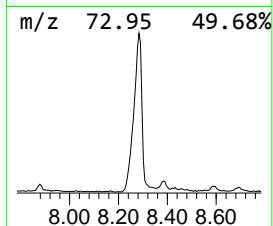
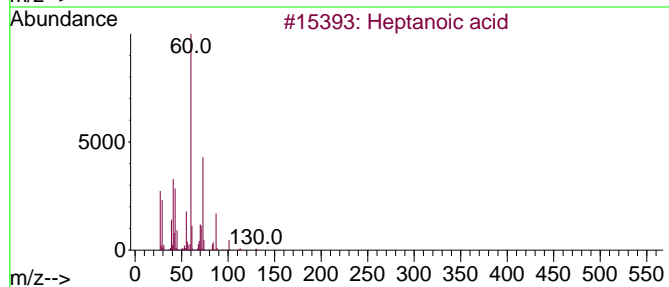
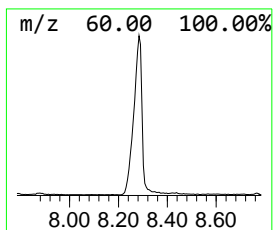
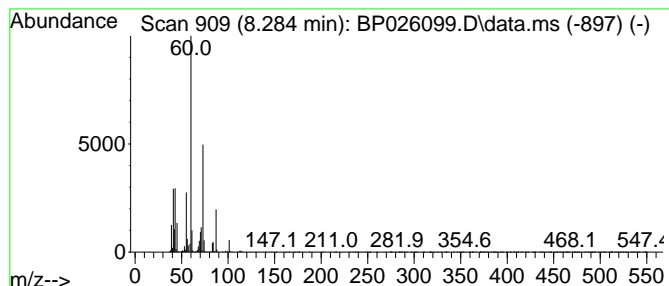
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 3 Heptanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.284	15.01 ng	1329440	1,4-Dichlorobenzene-d4	7.672

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Heptanoic acid	130	C7H14O2	000111-14-8	90
2		Butanoic acid, 3-methyl-	102	C5H10O2	000503-74-2	68
3		Hexanoic acid	116	C6H12O2	000142-62-1	64
4		Pentanoic acid	102	C5H10O2	000109-52-4	64
5		Pentanoic acid, 3-methyl-	116	C6H12O2	000105-43-1	45



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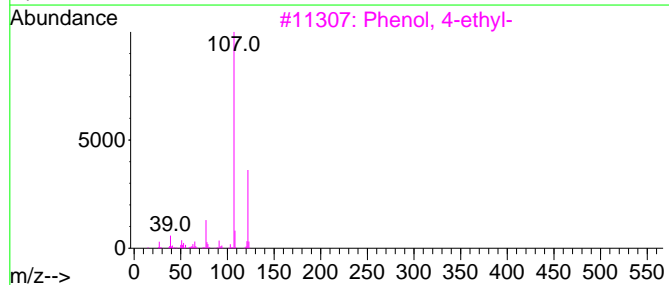
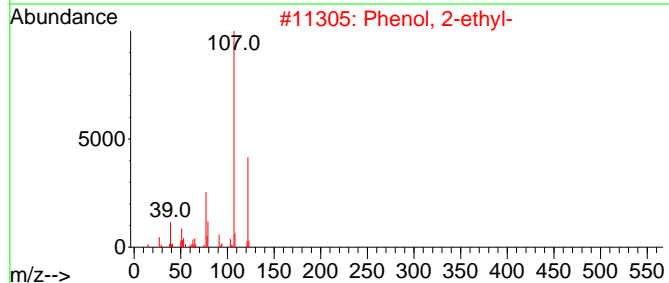
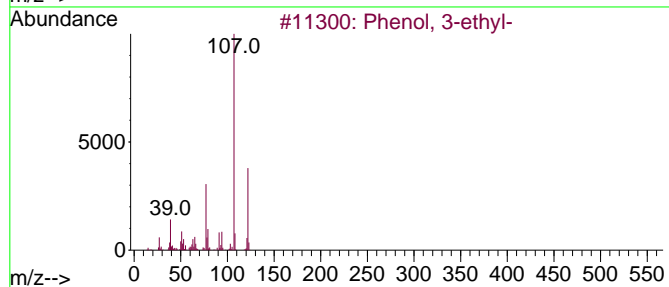
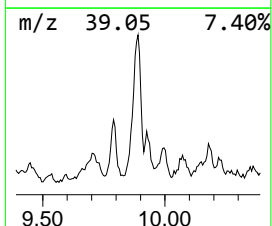
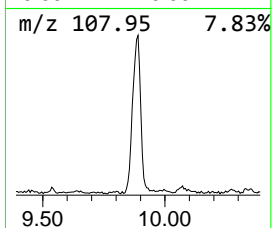
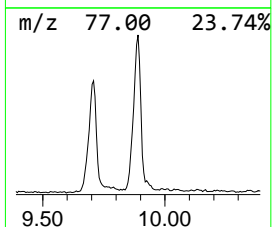
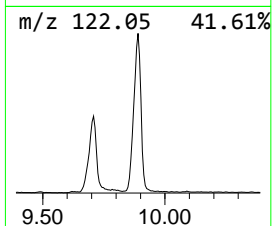
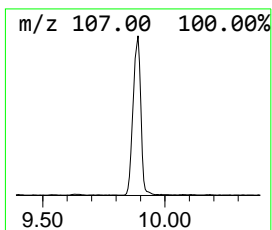
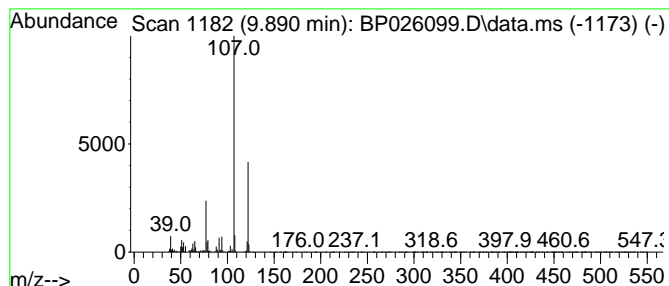
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 4 Phenol, 3-ethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.890	10.81 ng	1396750	Naphthalene-d8	10.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 3-ethyl-	122	C8H10O	000620-17-7	95
2			Phenol, 2-ethyl-	122	C8H10O	000090-00-6	91
3			Phenol, 4-ethyl-	122	C8H10O	000123-07-9	90
4			Phenol, 3,4-dimethyl-	122	C8H10O	000095-65-8	78
5			3,3-Dimethyl-6-methylenecyclohexene	122	C9H14	020185-16-4	72



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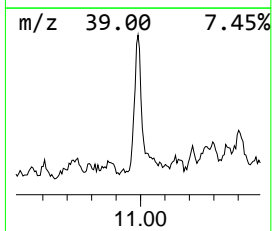
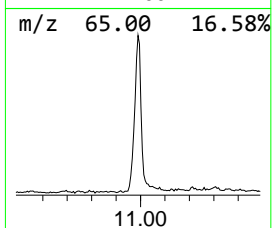
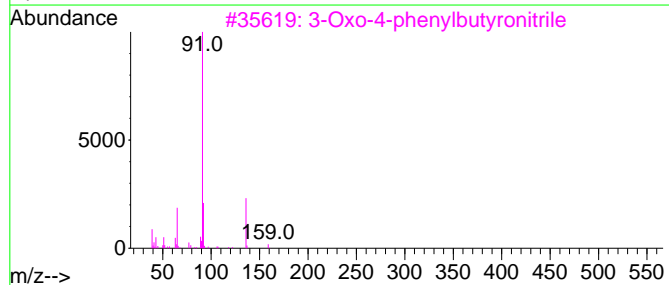
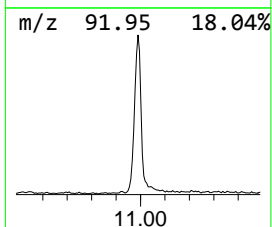
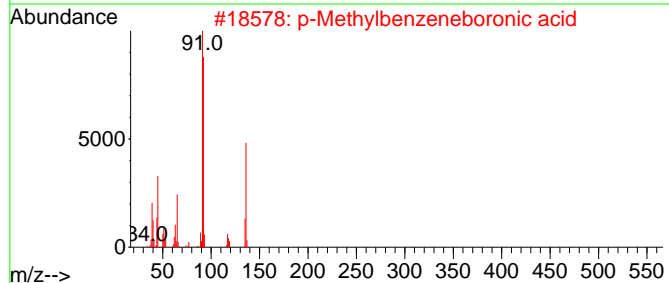
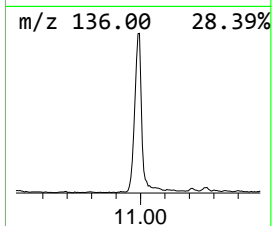
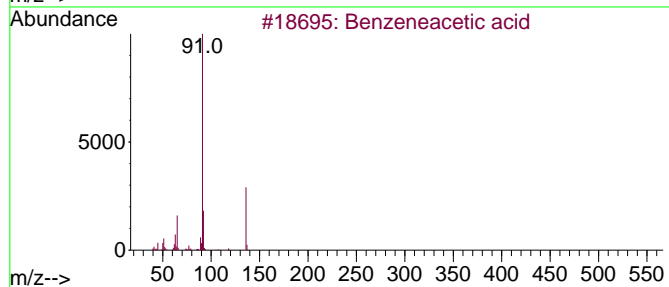
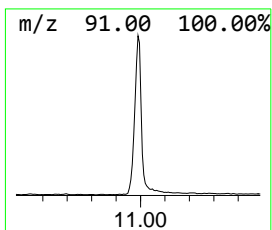
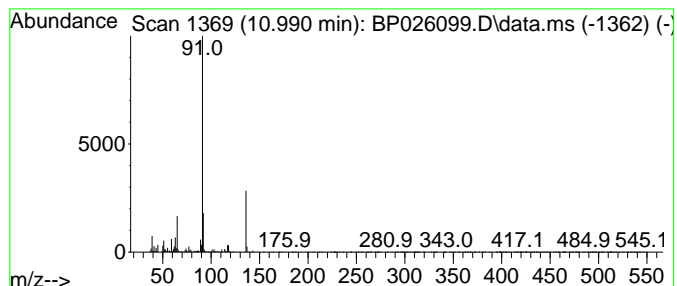
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TIC Library : C:\Database\NIST20.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 5 Benzeneacetic acid Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.990	8.90 ng	1149940	Naphthalene-d8	10.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid	136	C8H8O2	000103-82-2	93
2			p-Methylbenzeneboronic acid	136	C7H9B02	005720-05-8	52
3			3-Oxo-4-phenylbutyronitrile	159	C10H9N0	019212-27-2	43
4			4-Benzyloxybromobenzene	262	C13H11Br0	006793-92-6	43
5			Benzene, (iodomethyl)-	218	C7H7I	000620-05-3	43



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 ClientSampleId :
 SEEP-1

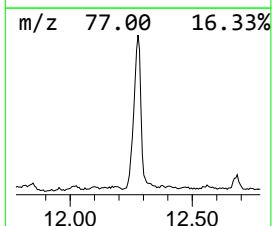
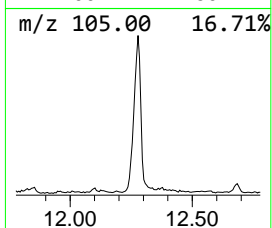
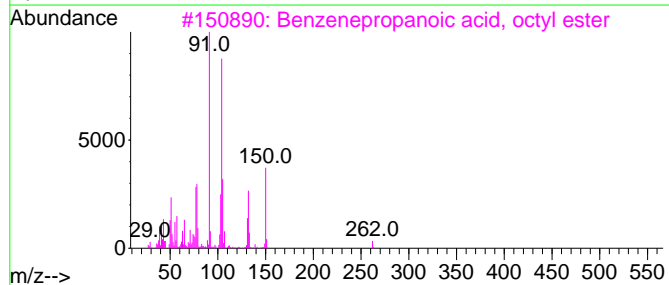
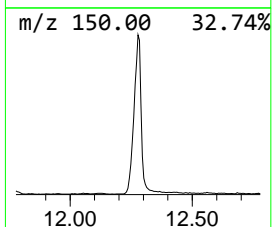
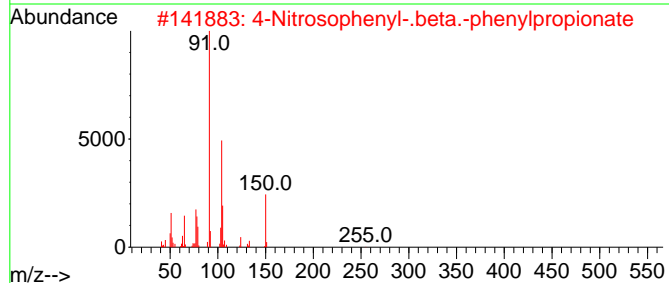
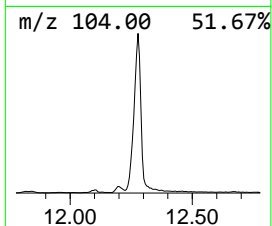
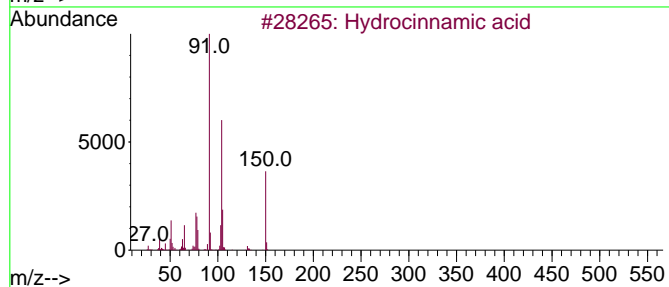
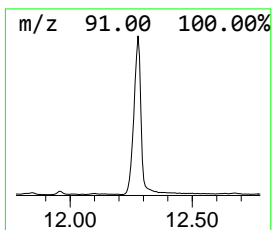
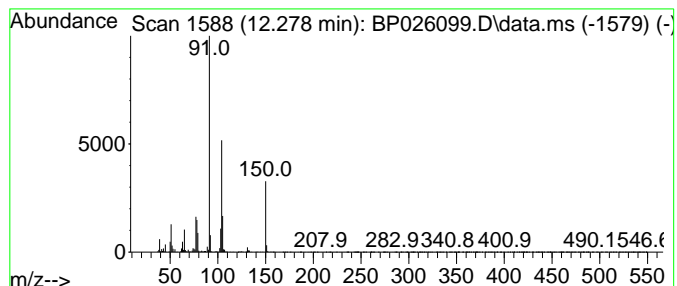
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 7 Hydrocinnamic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.278	17.08 ng	2206940	Naphthalene-d8	10.443

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hydrocinnamic acid	150	C9H10O2	000501-52-0	96
2		4-Nitrosophenyl-.beta.-phenylpro...	255	C15H13NO3	1000129-40-0	86
3		Benzenepropanoic acid, octyl ester	262	C17H26O2	037826-57-6	72
4		Benzenepropanoic acid, silver(1+...	256	C9H9AgO2	075112-79-7	53
5		2-Methylbenzyl alcohol, trifluor...	218	C10H9F3O2	1000364-57-0	43



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 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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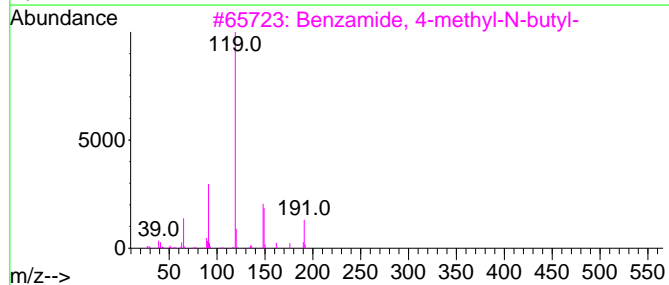
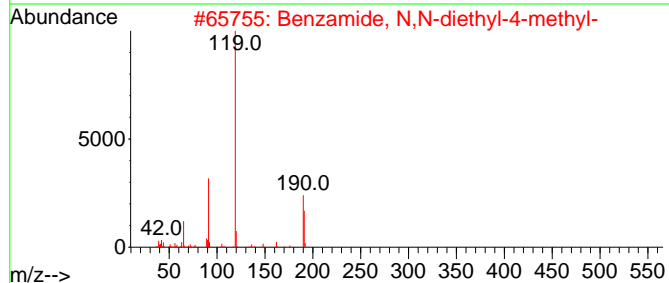
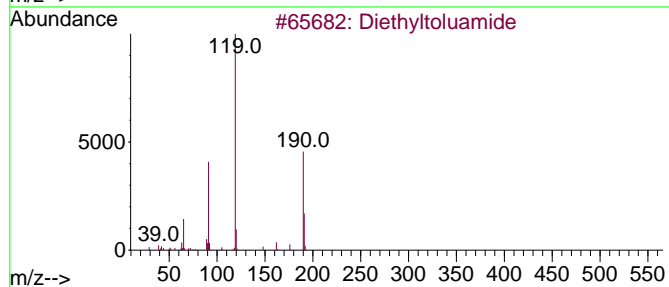
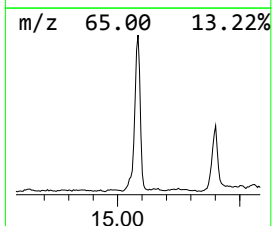
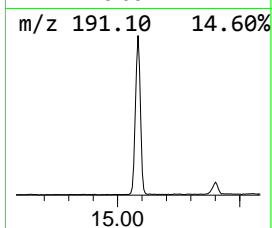
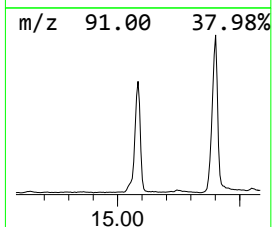
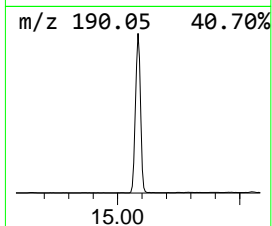
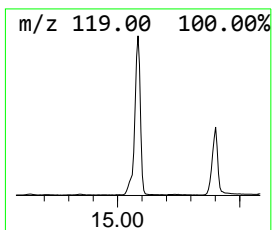
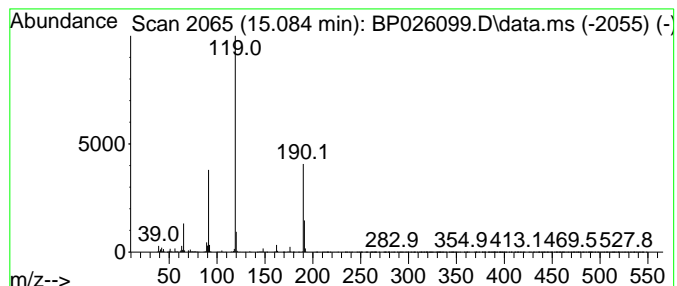
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 8 Diethyltoluamide Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.084	15.71 ng	3272620	Acenaphthene-d10	14.313

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Diethyltoluamide	191	C12H17NO	000134-62-3	96
2		Benzamide, N,N-diethyl-4-methyl-	191	C12H17NO	002728-05-4	93
3		Benzamide, 4-methyl-N-butyl-	191	C12H17NO	1000407-46-6	76
4		Benzamide, 3-methyl-N-butyl-	191	C12H17NO	1000407-41-1	76
5		Benzamide, 2-methyl-N-methyl-N-p...	191	C12H17NO	1000421-44-4	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
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 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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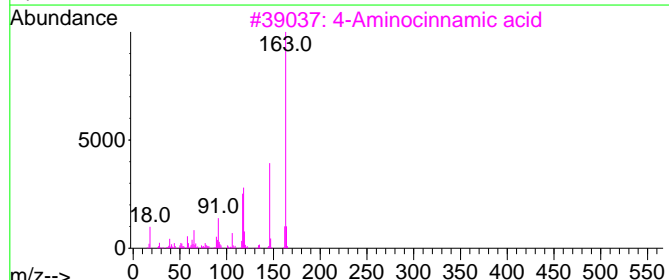
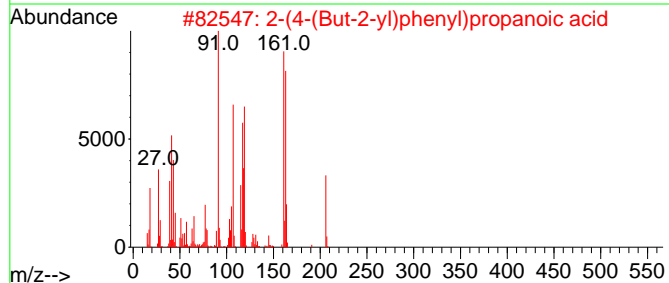
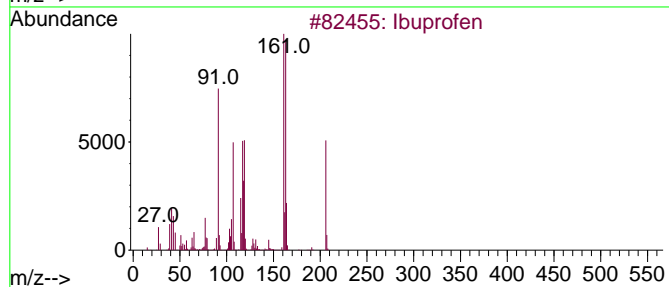
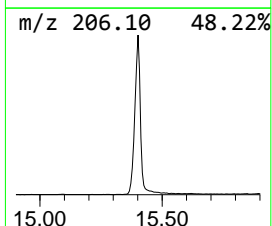
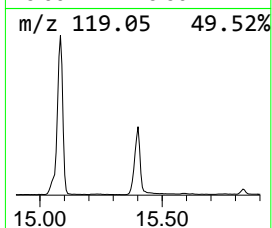
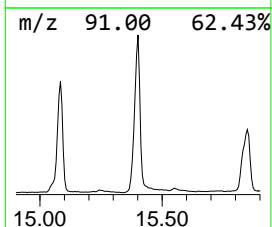
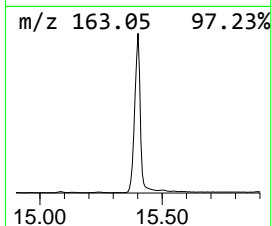
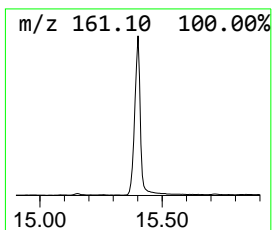
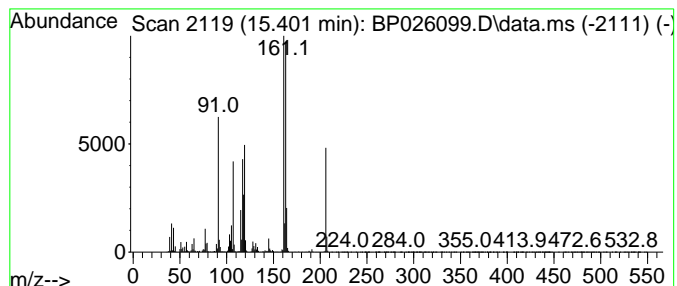
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 9 Ibuprofen Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.401	33.61 ng	6999860	Acenaphthene-d10	14.313

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Ibuprofen	206	C13H18O2	015687-27-1	99
2		2-(4-(But-2-yl)phenyl)propanoic ...	206	C13H18O2	064451-76-9	98
3		4-Aminocinnamic acid	163	C9H9NO2	002393-18-2	22
4		Thiocoumarin, 4,4,6-trimethyl-	206	C12H14O5	1000132-62-4	22
5		3-(3-Thienyl)pyridine	161	C9H7NS	021308-81-6	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
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Instrument :
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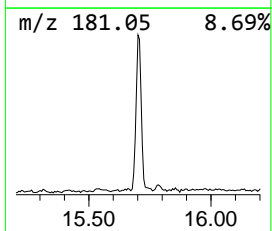
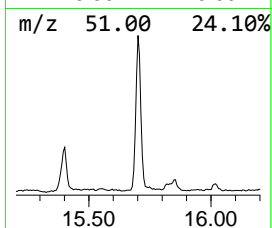
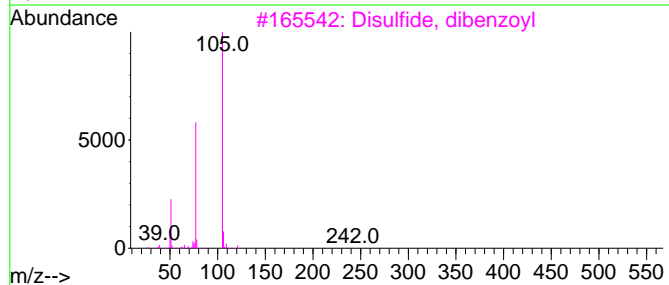
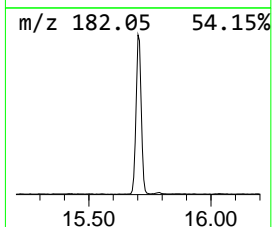
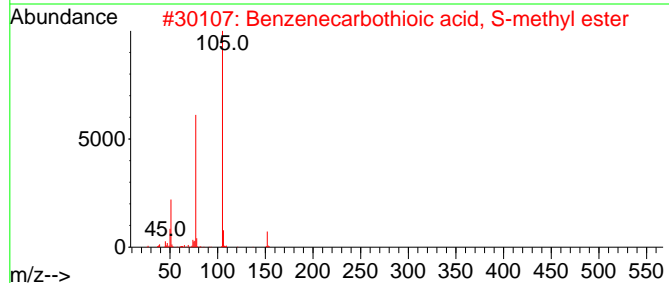
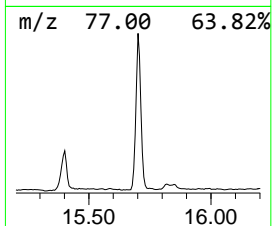
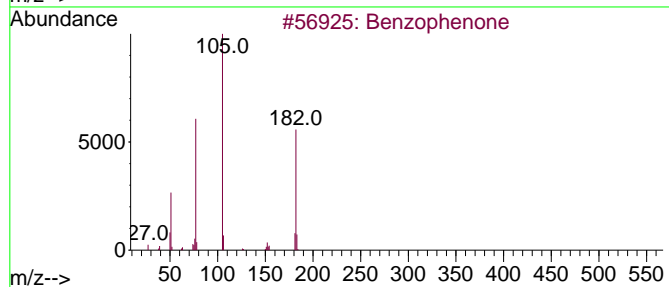
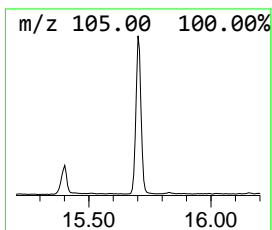
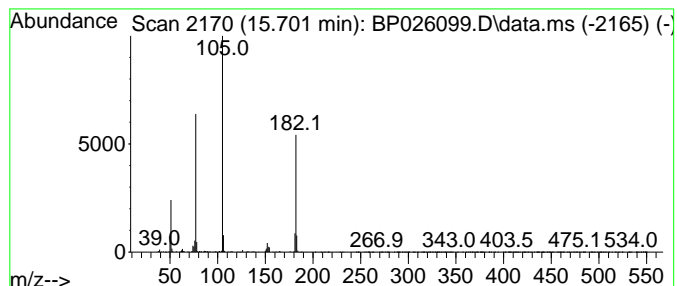
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TIC Library : C:\Database\NIST20.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 10 Benzophenone Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.701	9.10 ng	1894810	Acenaphthene-d10	14.313

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzophenone	182	C13H10O	000119-61-9	96
2		Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	50
3		Disulfide, dibenzoyl	274	C14H10O2S2	000644-32-6	47
4		acetic acid, 2-bromo-, 2-oxo-2-p...	256	C10H9BrO3	1000401-50-0	47
5		Benzenebutanoic acid, .gamma.-oxo-	178	C10H10O3	002051-95-8	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
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 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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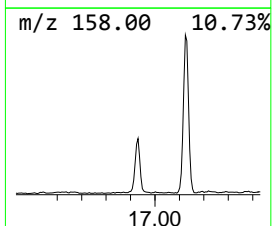
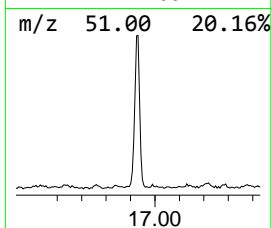
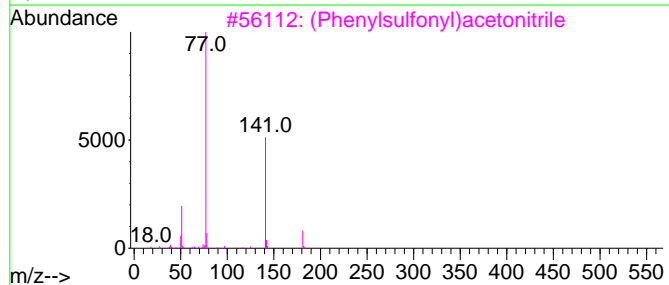
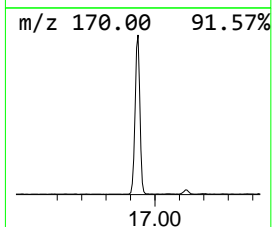
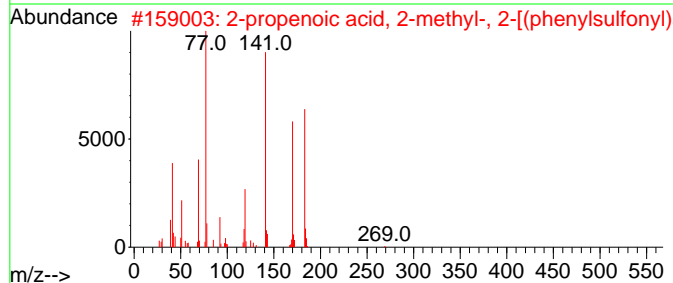
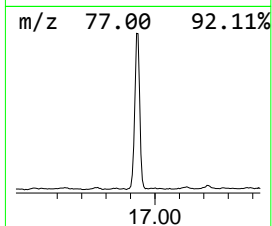
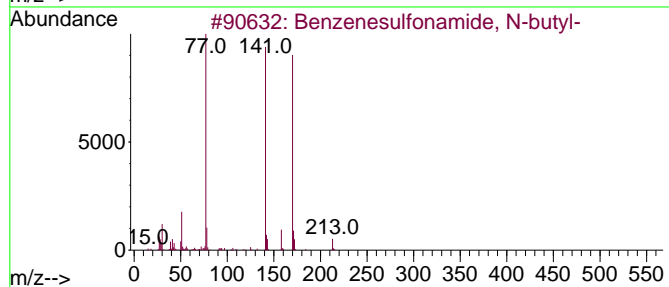
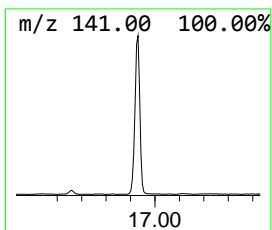
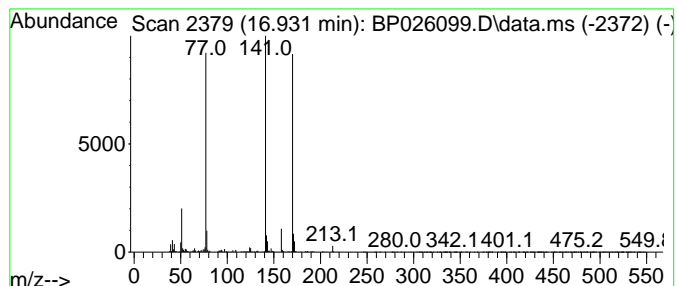
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 12 Benzenesulfonamide, N-butyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.931	8.59 ng	1892410	Phenanthrene-d10	17.131

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzenesulfonamide, N-butyl-	213	C10H15NO2S	003622-84-2	90
2		2-propenoic acid, 2-methyl-, 2-[...]	269	C12H15NO4S	1000401-71-8	64
3		(Phenylsulfonyl)acetonitrile	181	C8H7NO2S	007605-28-9	47
4		N-hexylbenzenesulfonamide	241	C12H19NO2S	007250-80-8	43
5		N-(2-Pyrazinyl)benzenesulfonamide	235	C10H9N3O2S	007471-20-7	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
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 Misc :
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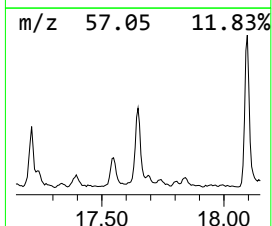
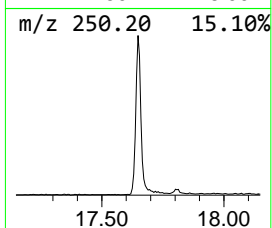
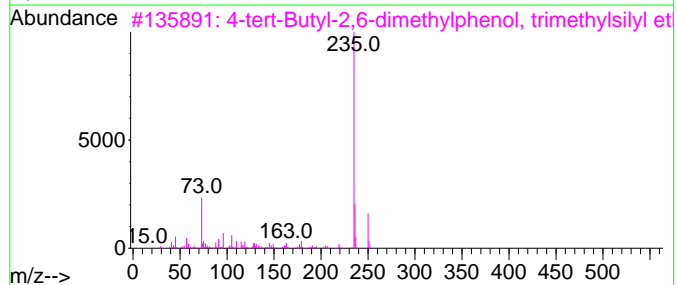
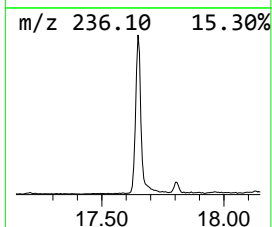
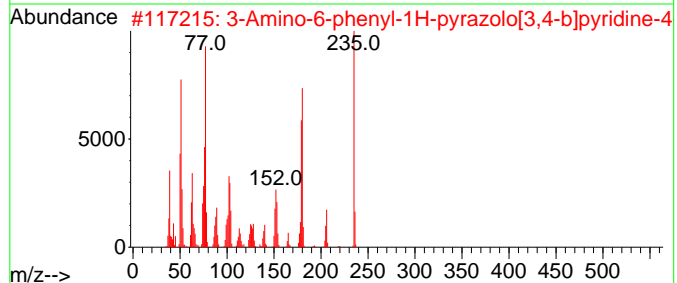
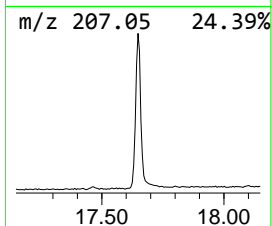
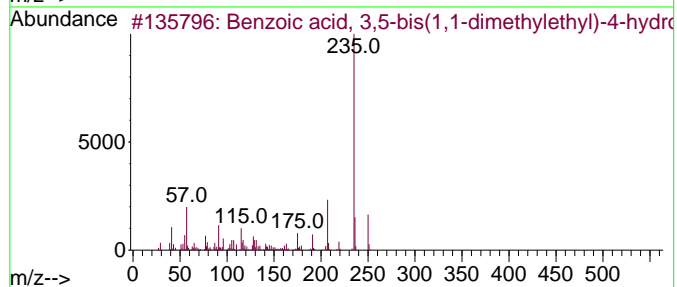
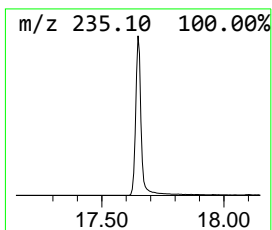
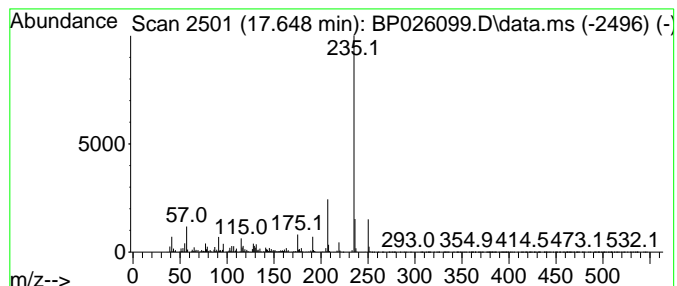
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TIC Library : C:\Database\NIST20.L
 TIC Integration Parameters: LSCINT.P

 Peak Number 13 Benzoic acid, 3,5-bis(1,1-d... Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.648	13.83 ng	3044800	Phenanthrene-d10	17.131

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 3,5-bis(1,1-dimeth...	250	C15H22O3	001421-49-4	96
2			3-Amino-6-phenyl-1H-pyrazolo[3,4...	235	C13H9N5	1000410-58-2	90
3			4-tert-Butyl-2,6-dimethylphenol,...	250	C15H26OSi	1000495-37-2	70
4			Hexobarbital-2(or 4)-methyl deri...	250	C13H18N2O3	1000123-51-5	59
5			6,7,8,9-Tetrahydro-1,2,3-trimeth...	250	C15H22O3	1000242-60-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
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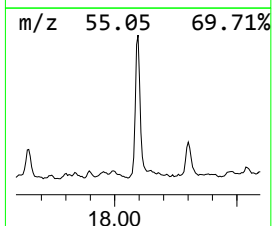
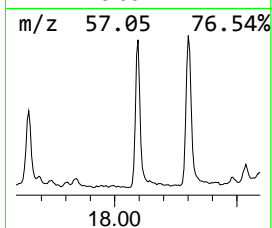
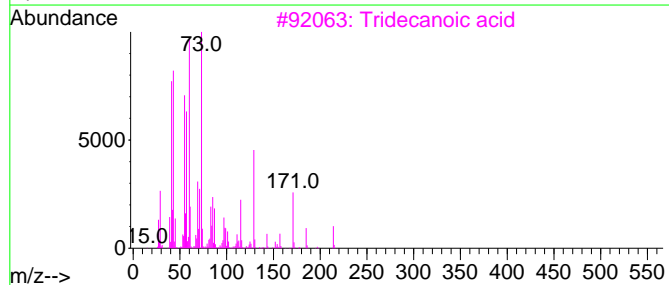
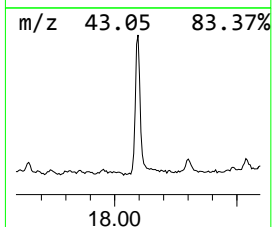
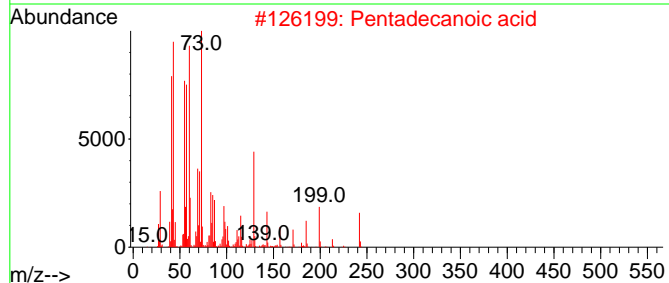
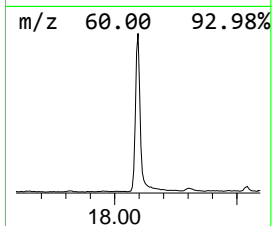
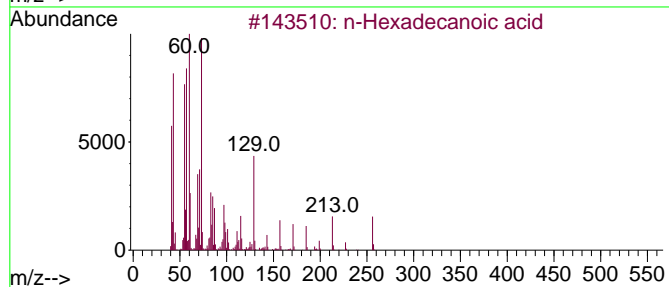
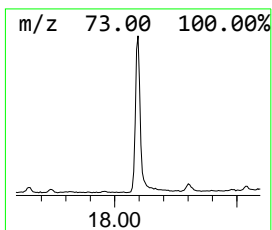
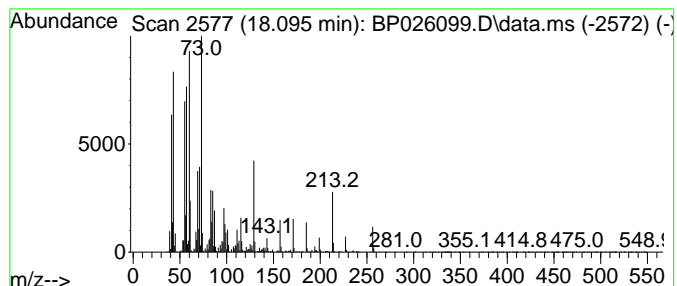
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 14 n-Hexadecanoic acid Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.095	13.49 ng	2970890	Phenanthrene-d10	17.131

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2		Pentadecanoic acid	242	C15H30O2	001002-84-2	93
3		Tridecanoic acid	214	C13H26O2	000638-53-9	93
4		Undecanoic acid	186	C11H22O2	000112-37-8	59
5		Tetradecanoic acid	228	C14H28O2	000544-63-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SEEP-1

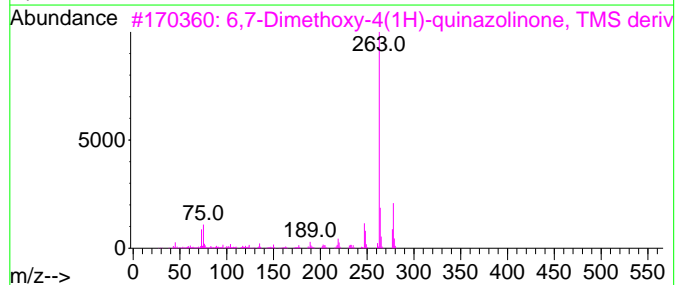
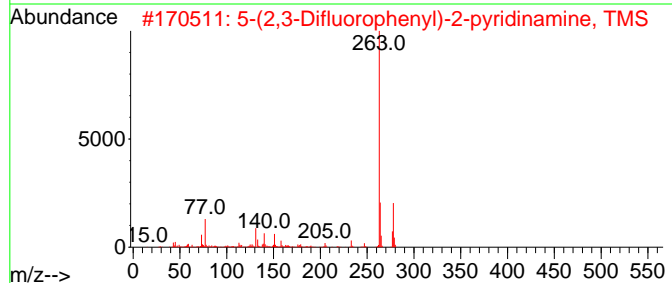
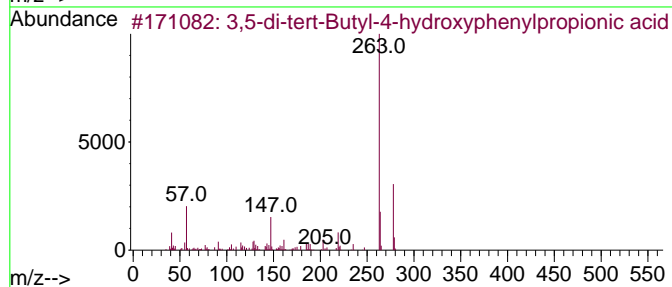
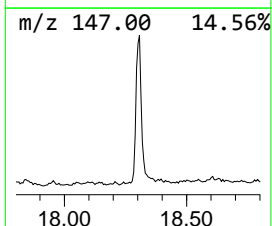
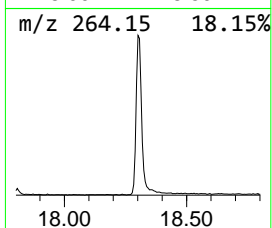
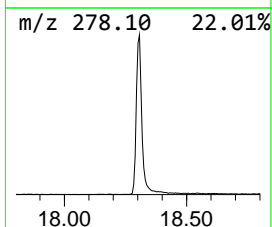
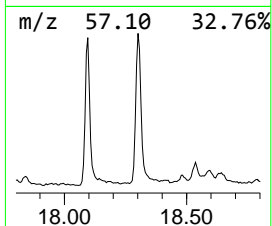
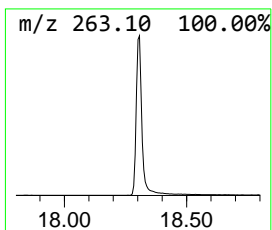
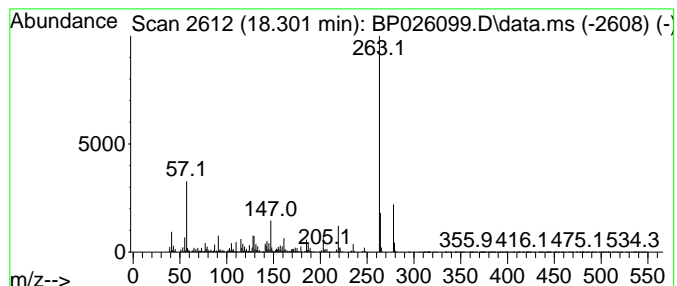
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 15 3,5-di-tert-Butyl-4-hydroxy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.301	14.05 ng	3093550	Phenanthrene-d10	17.131

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,5-di-tert-Butyl-4-hydroxypheny...	278	C17H26O3	020170-32-5	94
2		5-(2,3-Difluorophenyl)-2-pyridin...	278	C14H16F2N2Si	1000504-49-4	74
3		6,7-Dimethoxy-4(1H)-quinazolinon...	278	C13H18N2O3Si	1000479-65-5	72
4		4,6-Dinitro-1,1,3,3,5-pentamethy...	278	C14H18N2O4	000116-66-5	64
5		Cyanic acid, (1-methylethylidene...	278	C17H14N2O2	001156-51-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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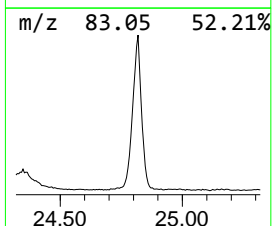
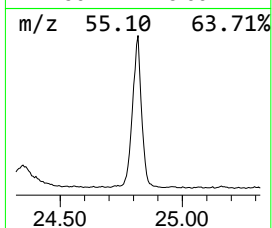
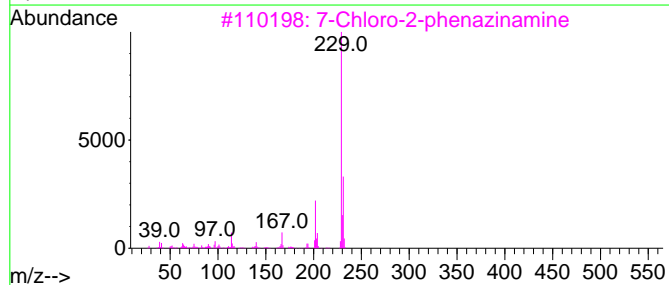
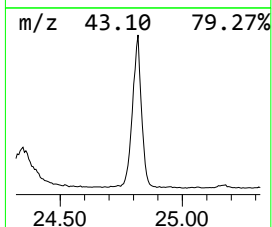
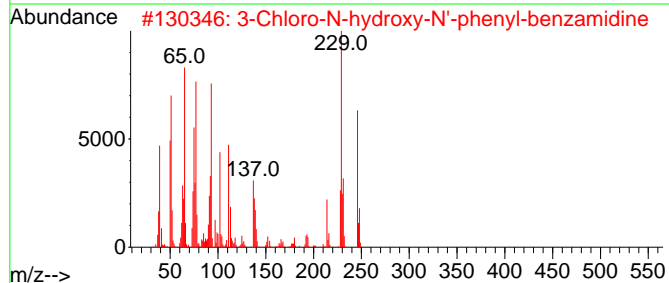
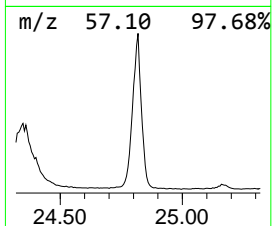
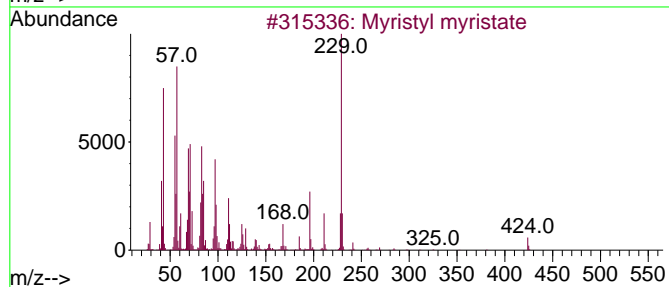
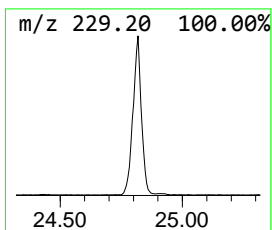
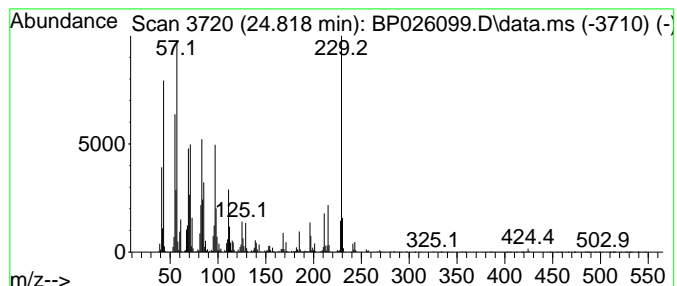
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 17 Myristyl myristate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.819	29.35 ng	8519050	Perylene-d12	24.913

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Myristyl myristate	424	C28H56O2	003234-85-3	93
2		3-Chloro-N-hydroxy-N'-phenyl-ben...	246	C13H11ClN2O	1000296-73-8	22
3		7-Chloro-2-phenazamine	229	C12H8ClN3	023677-11-4	22
4		10H-Phenothiazine, 3-methoxy-	229	C13H11NOS	001771-19-3	18
5		Glutaric acid, 2,2,3,3-tetraflu...	398	C19H30F4O4	1000394-01-1	16



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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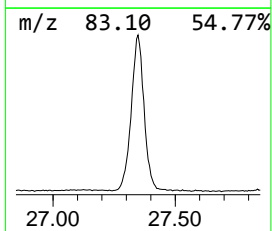
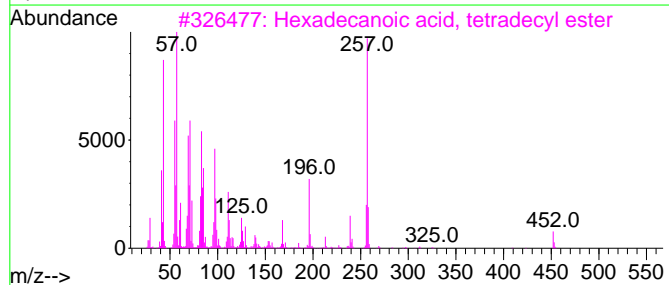
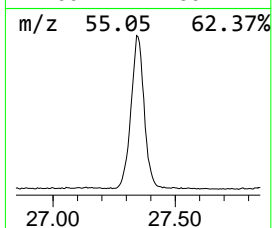
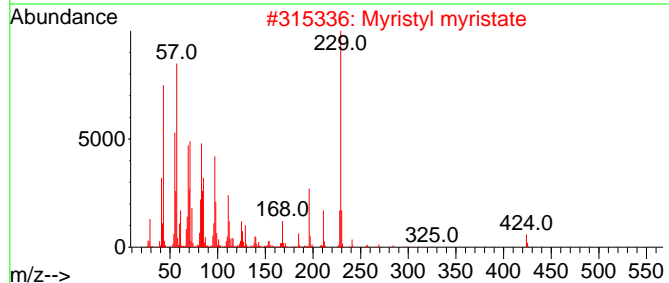
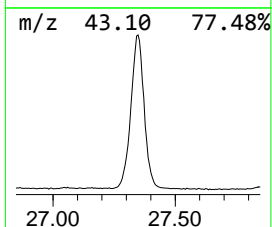
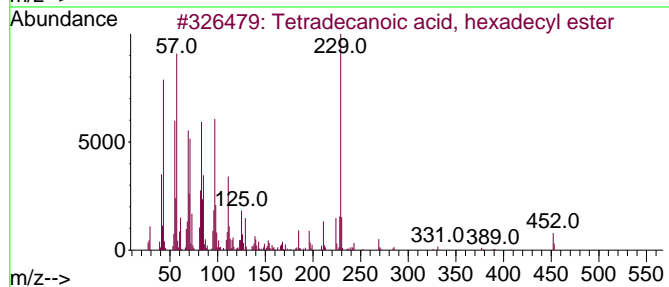
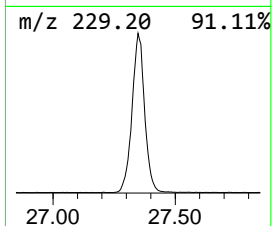
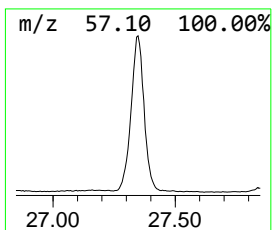
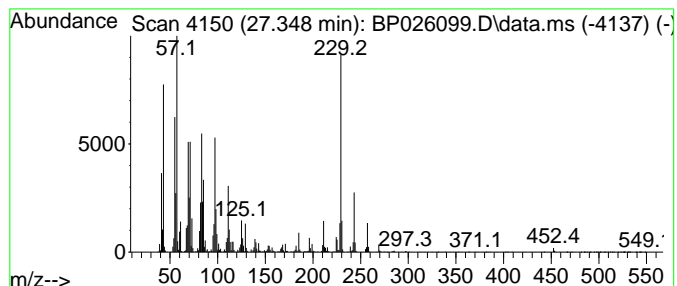
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 19 Tetradecanoic acid, hexadec... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
27.348	47.94 ng	13916300	Perylene-d12	24.913

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetradecanoic acid, hexadecyl ester	452	C30H60O2	002599-01-1	93
2		Myristyl myristate	424	C28H56O2	003234-85-3	87
3		Hexadecanoic acid, tetradecyl ester	452	C30H60O2	004536-26-9	38
4		6-Methyl-2-(methylthio)-4-pyrimi...	244	C9H16N2S2Si	1000479-10-2	18
5		10H-Phenothiazine, 3-methoxy-	229	C13H11NOS	001771-19-3	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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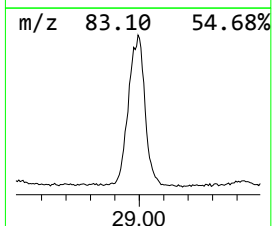
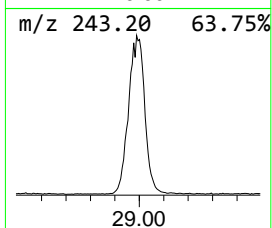
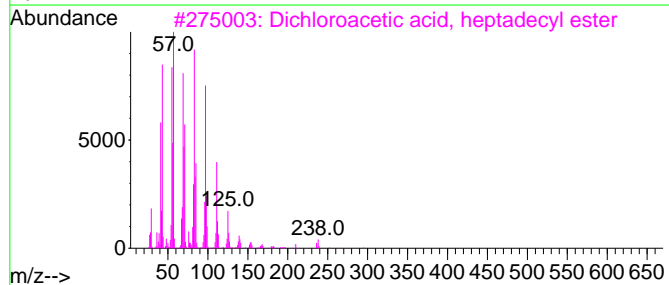
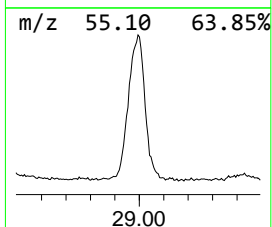
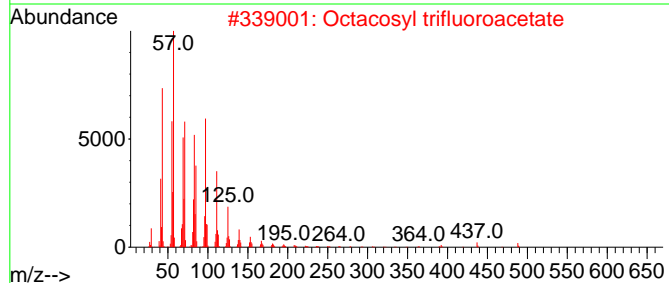
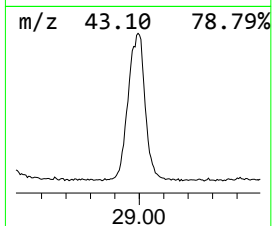
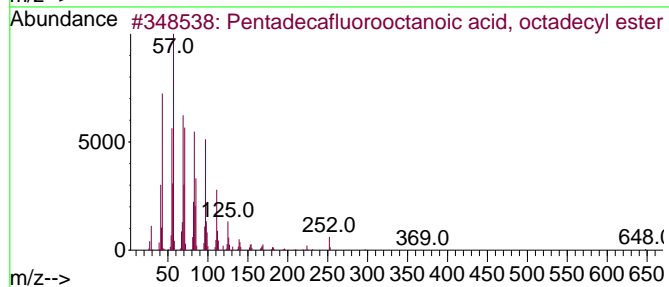
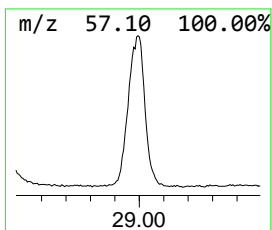
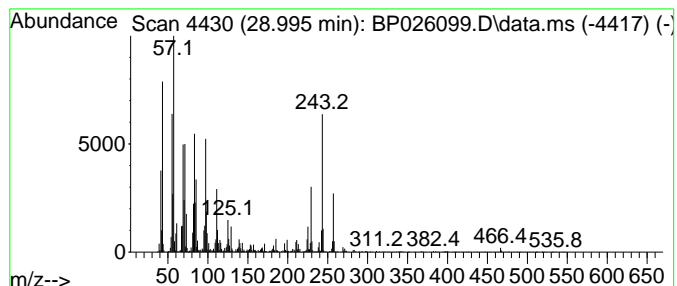
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 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

 Peak Number 20 Pentadecafluorooctanoic aci... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.995	23.09 ng	6704160	Perylene-d12	24.913

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	91
2		Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	90
3		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	90
4		Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	87
5		Hexadecyl pentadecanoate	466	C31H62O2	036617-33-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026099.D
 Acq On : 11 Nov 2025 20:59
 Operator : RC/JU
 Sample : Q3584-07
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Heptanoic acid	8.284	15.0	ng	1329440	1	7.672	1771080	20.0
Phenol, 3-ethyl-	9.890	10.8	ng	1396750	2	10.443	2584240	20.0
Benzeneacetic acid	10.990	8.9	ng	1149940	2	10.443	2584240	20.0
Hydrocinnamic acid	12.278	17.1	ng	2206940	2	10.443	2584240	20.0
Diethyltoluamide	15.084	15.7	ng	3272620	3	14.313	4165030	20.0
Ibuprofen	15.401	33.6	ng	6999860	3	14.313	4165030	20.0
Benzophenone	15.701	9.1	ng	1894810	3	14.313	4165030	20.0
Benzenesulfonam...	16.931	8.6	ng	1892410	4	17.131	4404490	20.0
Benzoic acid, 3...	17.648	13.8	ng	3044800	4	17.131	4404490	20.0
n-Hexadecanoic ...	18.095	13.5	ng	2970890	4	17.131	4404490	20.0
3,5-di-tert-But...	18.301	14.1	ng	3093550	4	17.131	4404490	20.0
Myristyl myristate	24.819	29.4	ng	8519050	6	24.913	5806110	20.0
Tetradecanoic a...	27.348	47.9	ng	13916300	6	24.913	5806110	20.0
Pentadecafluoro...	28.995	23.1	ng	6704160	6	24.913	5806110	20.0