

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
 Data File : BP026096.D
 Acq On : 11 Nov 2025 18:55
 Operator : RC/JU
 Sample : Q3569-02
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 MW-12

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_P\Methods\8270E-BP102925.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP026096.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.231	45	50	66	rVB	1559908	2145026	5.55%	1.009%
2	3.837	148	153	164	rBV	428664	665290	1.72%	0.313%
3	4.361	237	242	257	rVB3	219019	415946	1.08%	0.196%
4	5.043	350	358	365	rBV	323369	515723	1.34%	0.243%
5	5.155	370	377	386	rVB2	337667	571141	1.48%	0.269%
6	5.296	394	401	410	rBV	3264258	4794656	12.41%	2.255%
7	5.960	508	514	523	rBV2	202245	402554	1.04%	0.189%
8	6.737	626	646	652	rBV4	117349	466645	1.21%	0.219%
9	6.849	658	665	677	rBV	2023430	3372468	8.73%	1.586%
10	7.225	720	729	737	rBV	1118309	1995545	5.17%	0.938%
11	7.678	796	806	810	rBV	1398976	2540083	6.58%	1.194%
12	8.466	934	940	946	rBV4	325205	662214	1.71%	0.311%
13	8.649	964	971	975	rBV2	340032	650327	1.68%	0.306%
14	8.766	985	991	992	rBV2	300213	450306	1.17%	0.212%
15	8.807	992	998	1004	rVB	3193884	5685539	14.72%	2.674%
16	9.296	1072	1081	1087	rBV	674232	1356516	3.51%	0.638%
17	9.401	1087	1099	1105	rBV2	1721586	3354777	8.68%	1.578%
18	9.860	1169	1177	1184	rVB3	605295	1149014	2.97%	0.540%
19	10.007	1194	1202	1208	rVV2	703245	1382565	3.58%	0.650%
20	10.331	1245	1257	1270	rVV3	377338	1396639	3.62%	0.657%
21	10.443	1270	1276	1282	rVV	1620486	2726104	7.06%	1.282%
22	10.607	1291	1304	1311	rBV5	237603	661969	1.71%	0.311%
23	11.484	1449	1453	1461	rVB	242948	480794	1.24%	0.226%
24	11.795	1497	1506	1512	rBV	3153231	5103877	13.21%	2.400%
25	11.966	1529	1535	1542	rBV5	271506	593208	1.54%	0.279%
26	12.919	1691	1697	1708	rVB	8902625	13997732	36.24%	6.582%
27	13.154	1733	1737	1743	rVB3	228246	415871	1.08%	0.196%
28	13.219	1743	1748	1752	rBV	292984	513204	1.33%	0.241%
29	13.542	1798	1803	1805	rBV2	543714	799624	2.07%	0.376%
30	13.572	1806	1808	1821	rVB2	555896	898461	2.33%	0.422%
31	13.701	1826	1830	1834	rVB	567454	777803	2.01%	0.366%
32	13.872	1853	1859	1864	rBV	1377544	2750983	7.12%	1.294%
33	14.142	1899	1905	1912	rBV	936665	1628455	4.22%	0.766%
34	14.237	1917	1921	1927	rVV2	465827	910420	2.36%	0.428%
35	14.313	1928	1934	1940	rVB	2589032	3764105	9.74%	1.770%
36	14.595	1978	1982	1986	rVB	467940	586813	1.52%	0.276%

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37	14.660	1988	1993	1999	rVV2	769351	1330910	3.45%	0.626%
38	14.742	2003	2007	2012	rVB4	250574	432600	1.12%	0.203%
39	14.936	2037	2040	2045	rVB	519756	654580	1.69%	0.308%
40	15.089	2062	2066	2070	rVB	485381	617094	1.60%	0.290%
41	15.401	2113	2119	2125	rVB	1831237	2725657	7.06%	1.282%
42	15.601	2147	2153	2166	rBV2	665107	1656979	4.29%	0.779%
43	15.707	2166	2171	2173	rVV	768280	1038744	2.69%	0.488%
44	15.731	2173	2175	2180	rVB	721825	875665	2.27%	0.412%
45	15.842	2184	2194	2204	rBV2	6788925	15256651	39.50%	7.174%
46	16.101	2221	2238	2249	rVB	10438836	38628580	100.00%	18.164%
47	16.395	2282	2288	2297	rBV	1325810	2331935	6.04%	1.097%
48	16.766	2346	2351	2358	rVB2	923435	1511442	3.91%	0.711%
49	16.930	2374	2379	2384	rBV4	274131	547885	1.42%	0.258%
50	17.130	2408	2413	2419	rVB2	2495365	4135756	10.71%	1.945%
51	17.283	2435	2439	2444	rVB2	283875	408875	1.06%	0.192%
52	17.648	2497	2501	2503	rBV	478789	615669	1.59%	0.290%
53	17.695	2503	2509	2518	rVV2	5939049	9614068	24.89%	4.521%
54	17.777	2518	2523	2529	rVB	2781394	3923844	10.16%	1.845%
55	18.095	2572	2577	2585	rVB2	868957	1339111	3.47%	0.630%
56	18.313	2611	2614	2621	rBV2	287022	466329	1.21%	0.219%
57	18.860	2704	2707	2715	rVB3	366087	652830	1.69%	0.307%
58	18.954	2720	2723	2731	rVB4	267629	480905	1.24%	0.226%
59	19.272	2773	2777	2783	rVB2	437122	672855	1.74%	0.316%
60	19.360	2788	2792	2799	rVB	1589904	2198676	5.69%	1.034%
61	19.683	2841	2847	2860	rVB	9441002	13490604	34.92%	6.344%
62	19.872	2873	2879	2886	rVB	13250487	21267050	55.06%	10.000%
63	20.066	2908	2912	2916	rBV2	327424	440528	1.14%	0.207%
64	20.207	2930	2936	2940	rVB	1109427	1682893	4.36%	0.791%
65	20.742	3023	3027	3033	rBV	1375598	1523939	3.95%	0.717%
66	21.277	3115	3118	3123	rVB	317801	410184	1.06%	0.193%
67	21.583	3164	3170	3180	rVB	3295855	5270285	13.64%	2.478%
68	24.912	3727	3736	3756	rVB	1817090	5876943	15.21%	2.764%

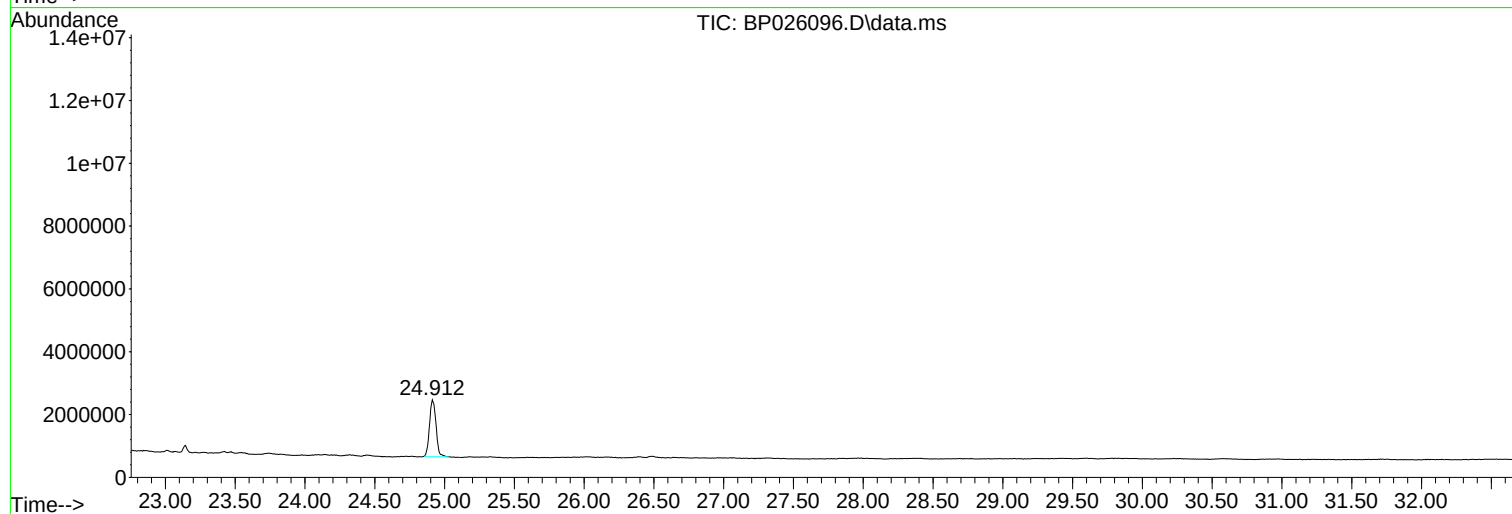
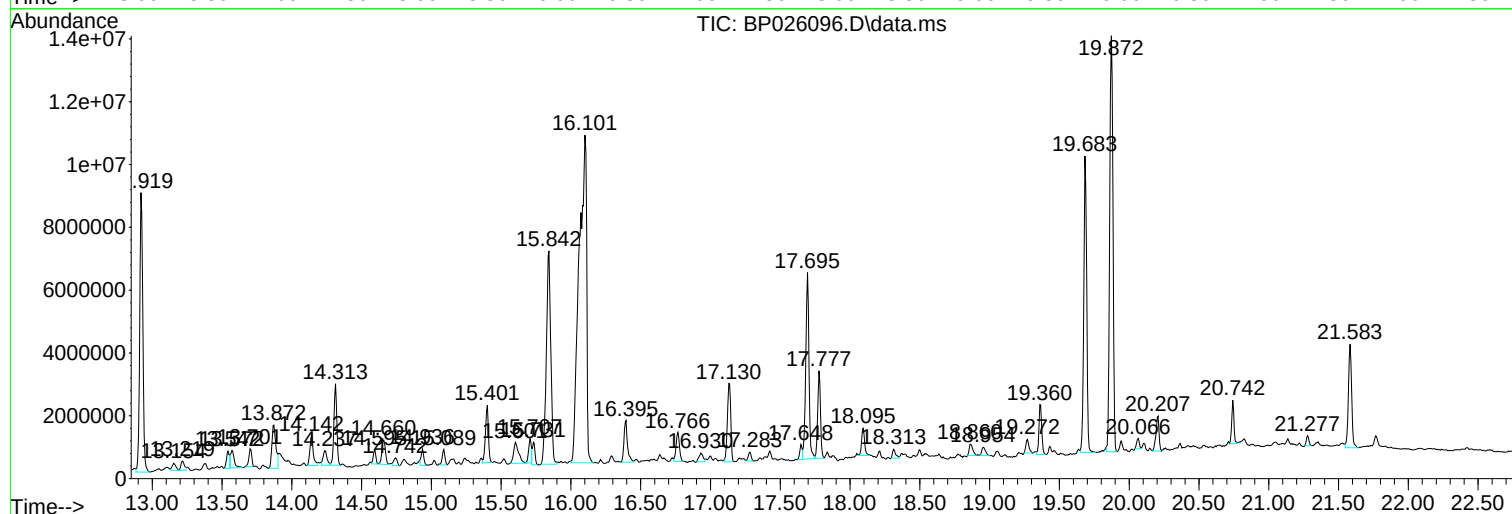
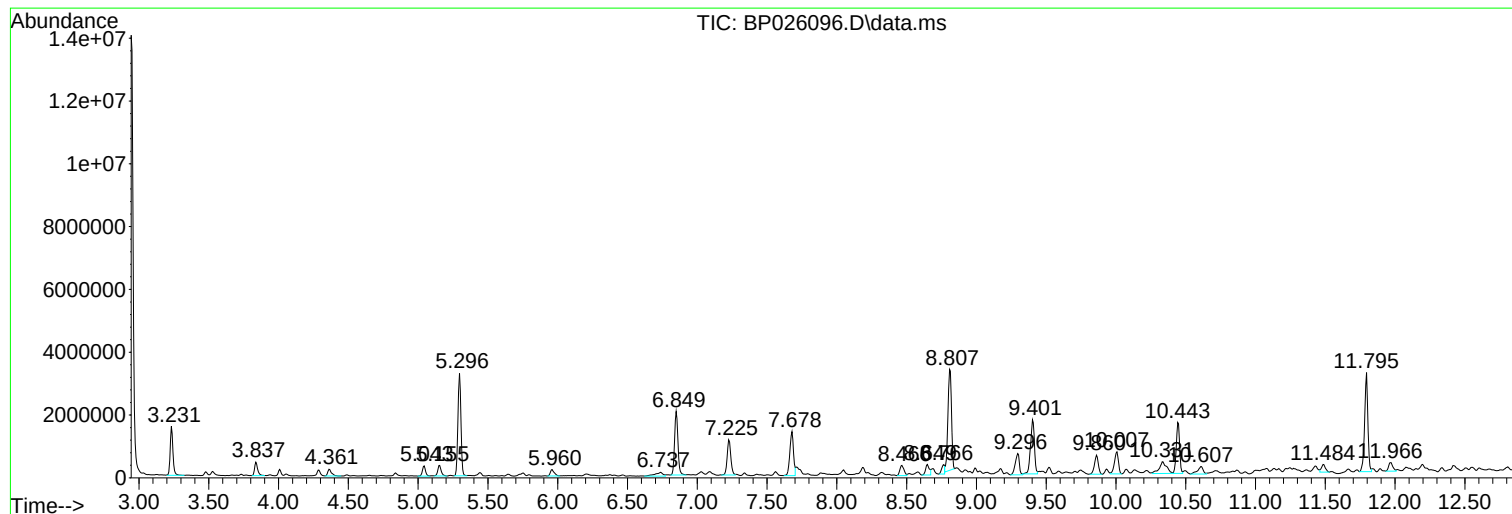
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Sample : Q3569-02
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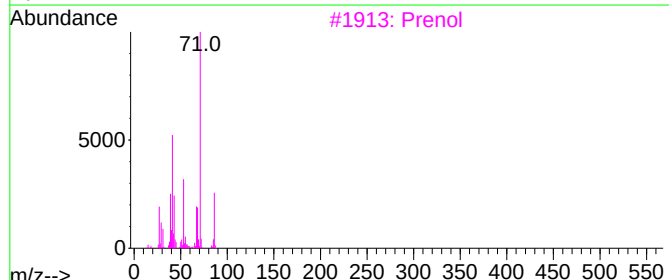
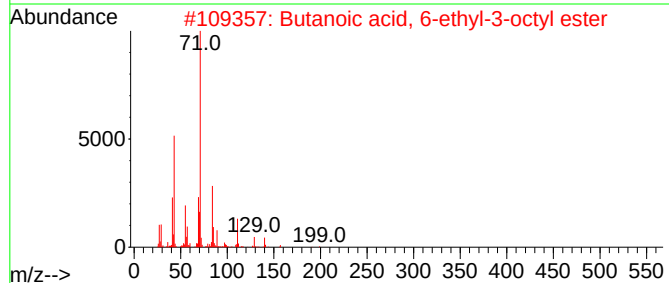
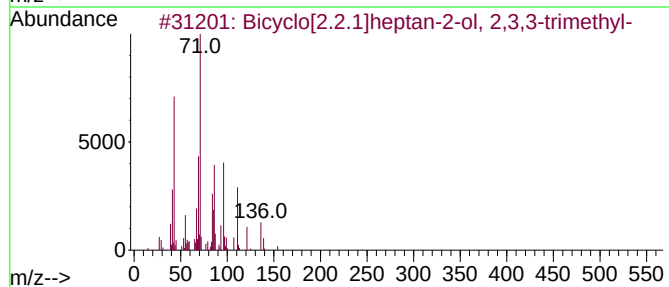
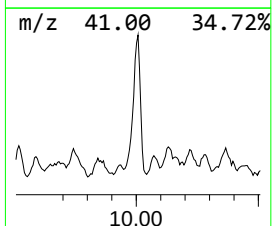
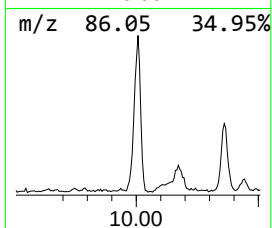
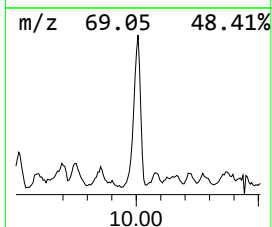
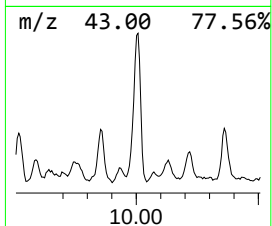
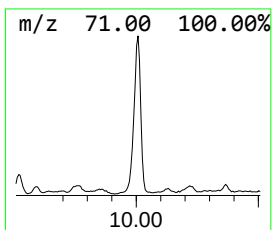
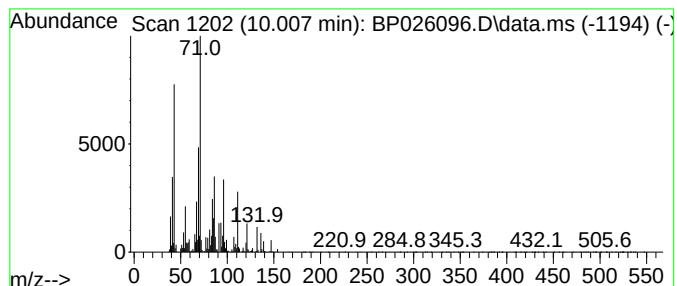
Instrument :
BNA_P
ClientSampleId :
MW-12

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 4 Bicyclo[2.2.1]heptan-2-ol, ... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.007	10.14 ng	1382570	Naphthalene-d8	10.443	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[2.2.1]heptan-2-ol, 2,3,3...	154	C10H18O	000465-31-6	96
2	Butanoic acid, 6-ethyl-3-octyl e...	228	C14H28O2	1000282-60-1	43
3	Prenol	86	C5H10O	000556-82-1	43
4	Z-1-Methoxy-2-hexene	114	C7H14O	1000279-60-9	38
5	Ether, methyl 1-tetradecenyl	226	C15H30O	026537-05-3	35



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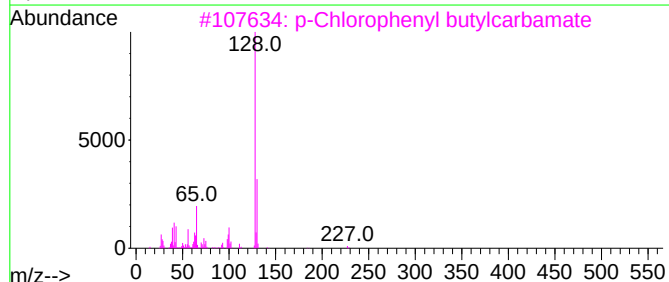
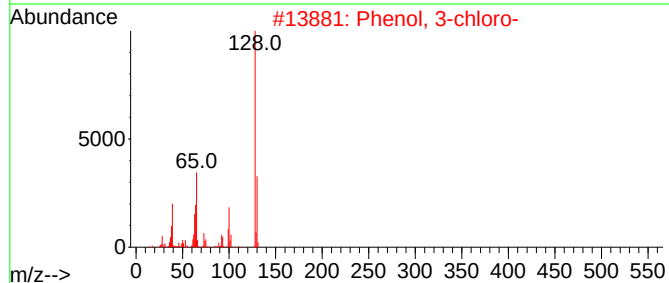
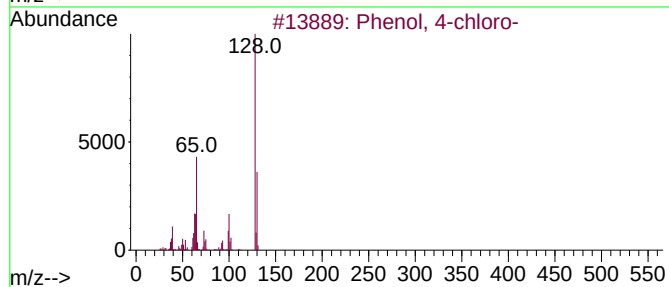
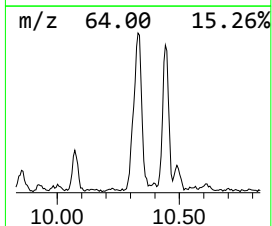
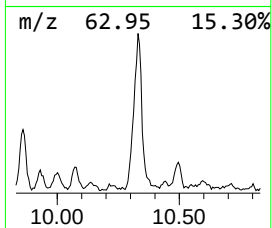
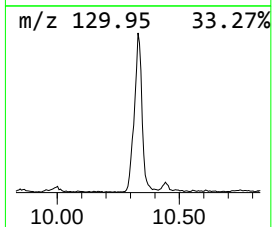
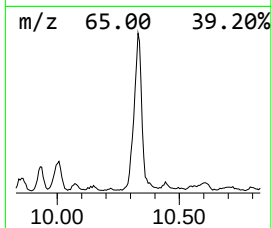
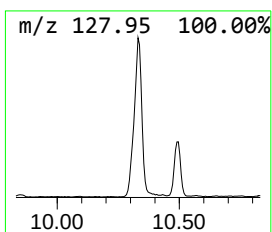
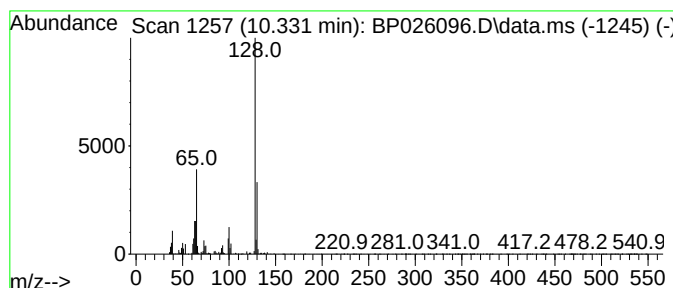
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Phenol, 4-chloro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.331	10.25 ng	1396640	Naphthalene-d8	10.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-chloro-	128	C6H5ClO	000106-48-9	97
2			Phenol, 3-chloro-	128	C6H5ClO	000108-43-0	94
3			p-Chlorophenyl butylcarbamate	227	C11H14ClNO2	132905-93-2	83
4			p-Chlorophenyl N-(cyanomethyl)ca...	210	C9H7ClN2O2	1000240-68-8	38
5			Phenol, 2-chloro-	128	C6H5ClO	000095-57-8	38



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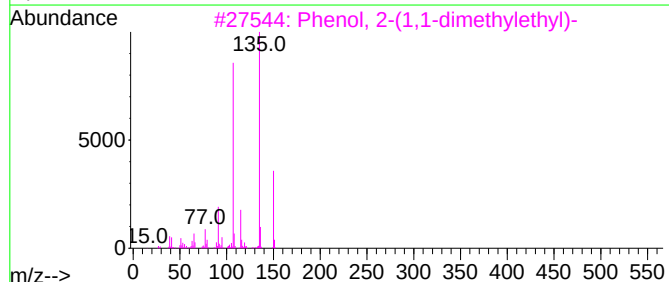
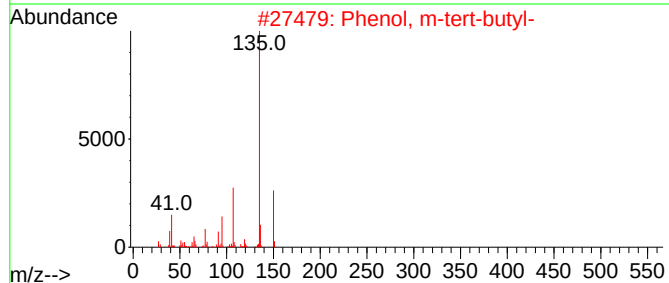
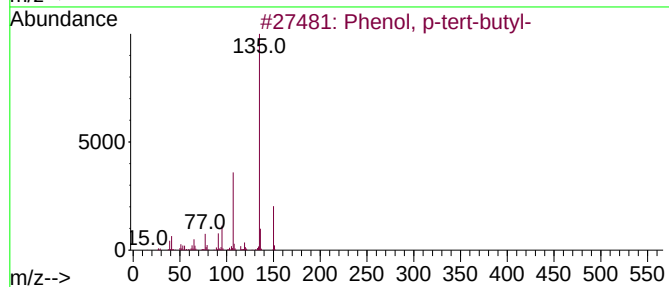
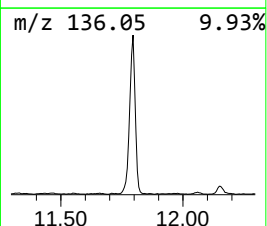
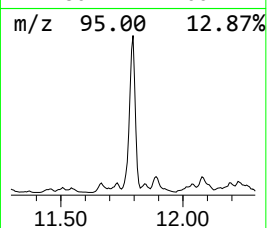
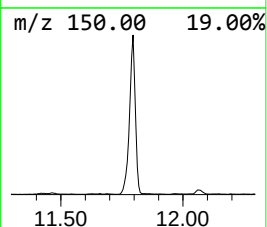
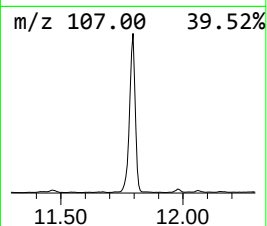
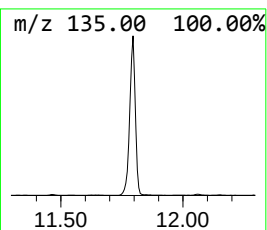
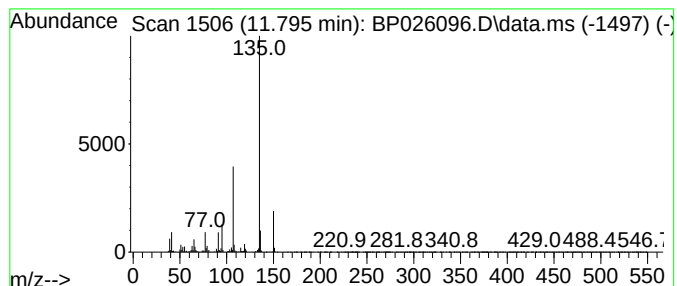
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 Phenol, p-tert-butyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.796	37.44 ng	5103880	Naphthalene-d8	10.443

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	97
2			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	94
3			Phenol, 2-(1,1-dimethylethyl)-	150	C10H14O	000088-18-6	74
4			Imidazo[1,2-c]pyrimidin-5-ol	135	C6H5N3O	055662-66-3	58
5			Phenol, 2-methyl-5-(1-methylethyl)-	150	C10H14O	000499-75-2	52



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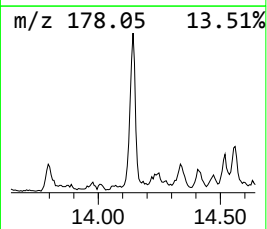
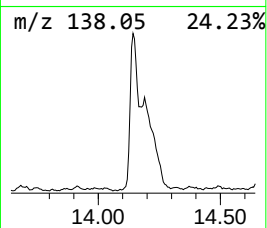
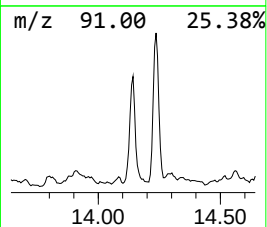
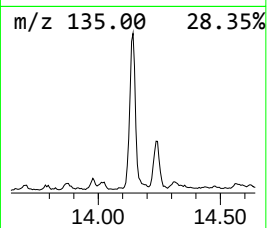
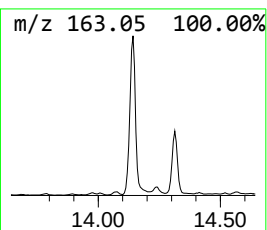
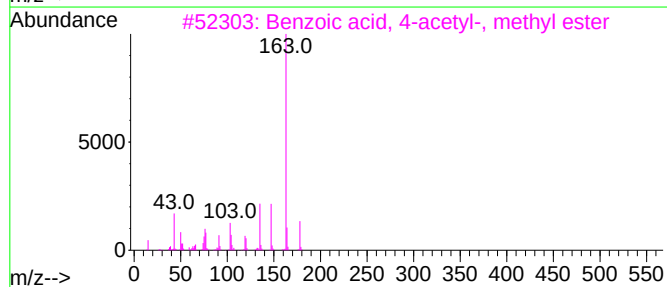
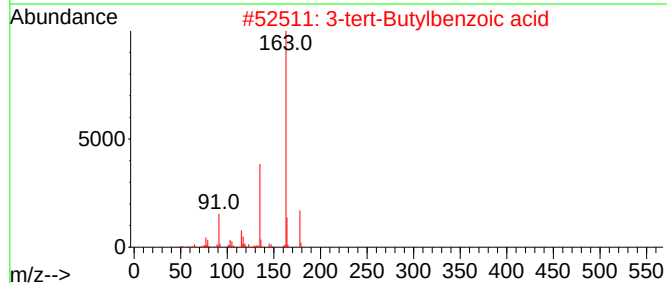
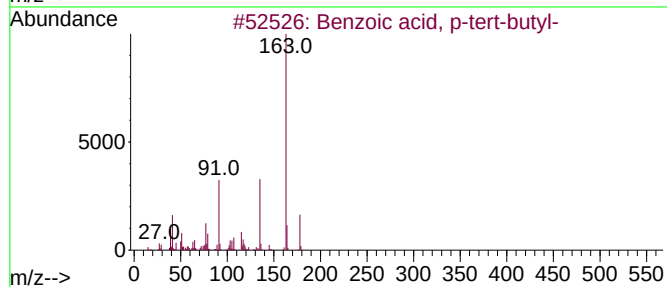
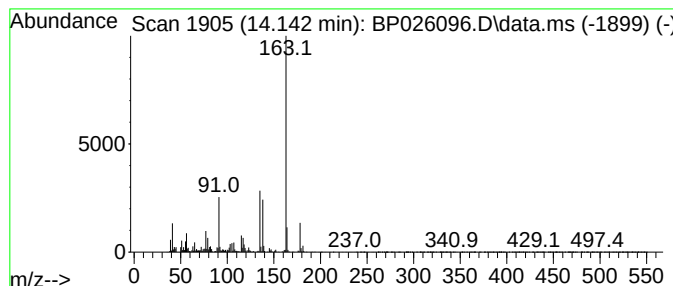
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Benzoic acid, p-tert-butyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.142	8.65 ng	1628460	Acenaphthene-d10	14.313	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzoic acid, p-tert-butyl-	178	C11H14O2	000098-73-7	95
2	3-tert-Butylbenzoic acid	178	C11H14O2	007498-54-6	76
3	Benzoic acid, 4-acetyl-, methyl ...	178	C10H10O3	003609-53-8	58
4	Terephthalaldehyde mono-(diethyl...	208	C12H16O3	081172-89-6	53
5	6-tert-Butyl-2,4-dimethylphenol	178	C12H18O	001879-09-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

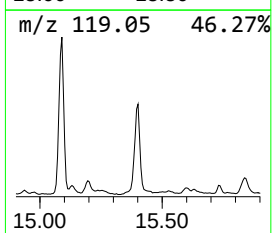
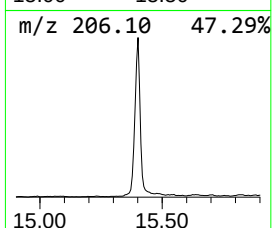
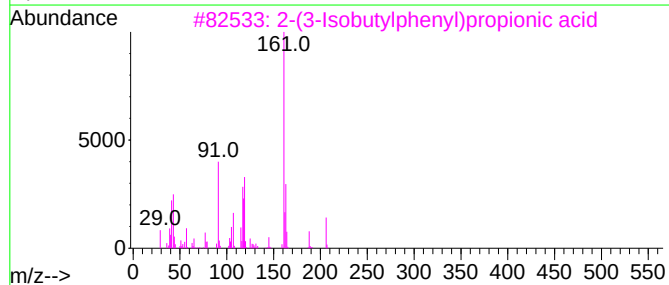
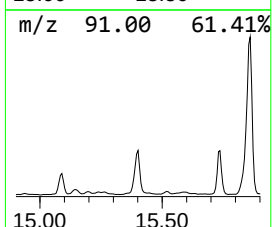
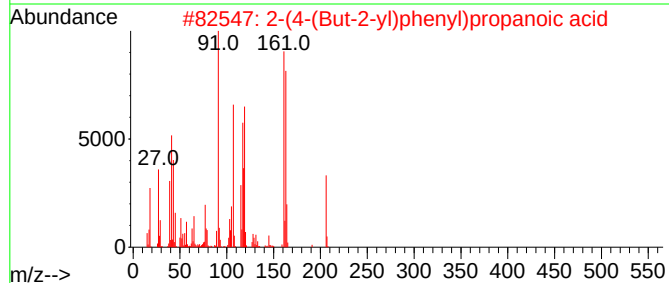
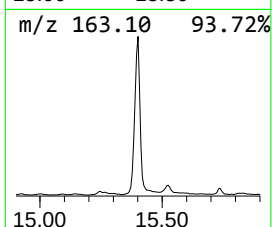
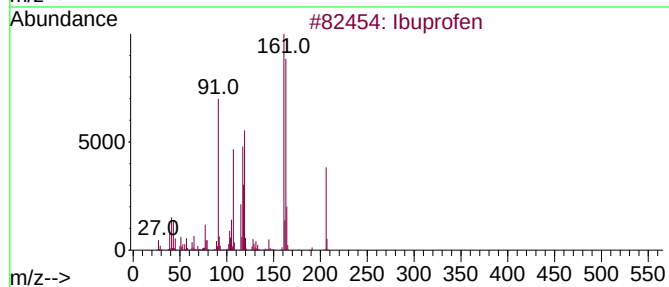
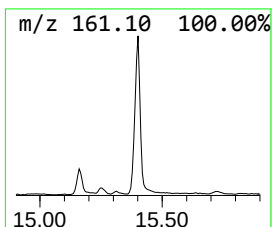
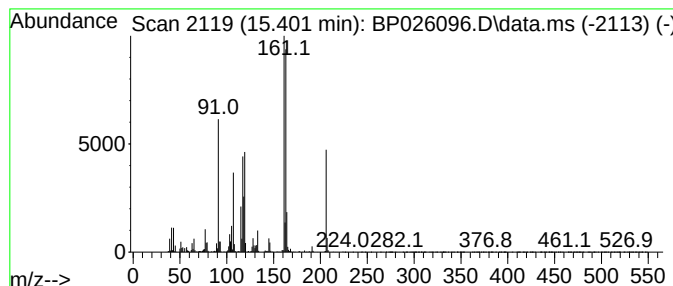
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Ibuprofen Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.401	14.48 ng	2725660	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	87
3	2-(3-Isobutylphenyl)propionic acid		206	C13H18O2	066622-47-7	50
4	Benzenamine, 2,4-dichloro-		161	C6H5Cl2N	000554-00-7	38
5	Benzenamine, 2,5-dichloro-		161	C6H5Cl2N	000095-82-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-12

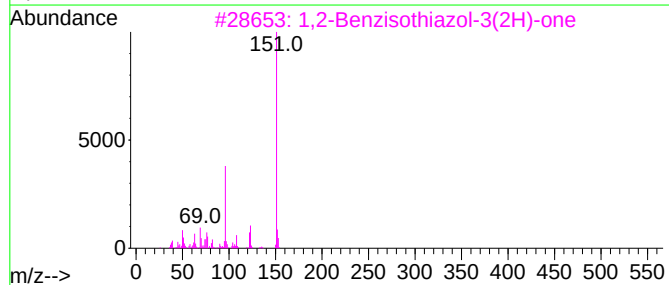
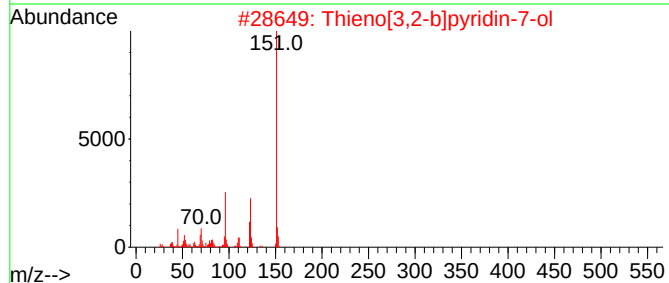
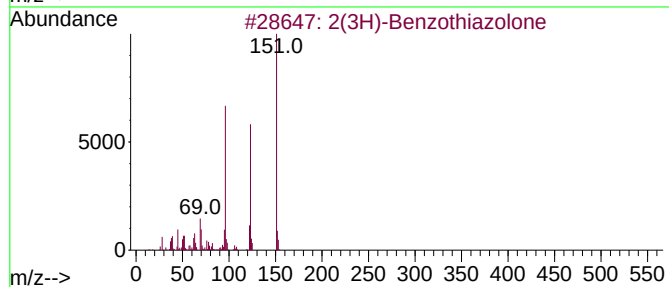
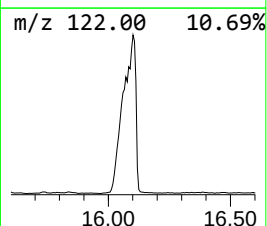
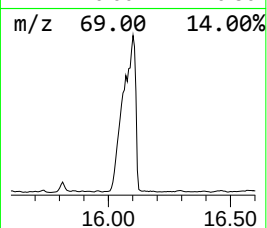
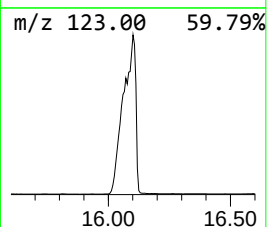
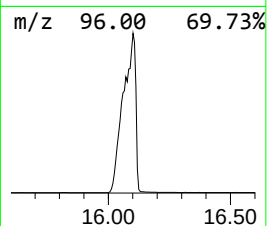
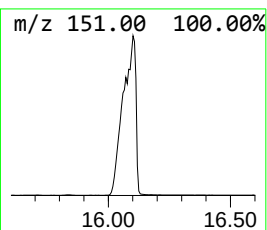
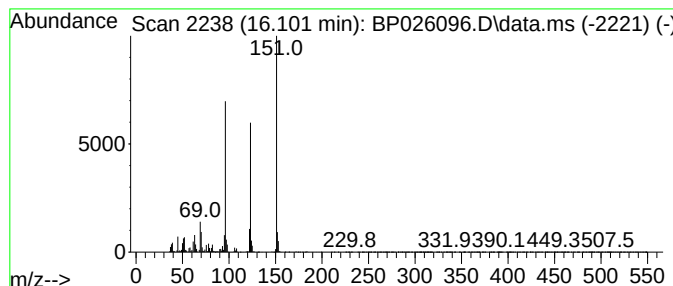
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 2(3H)-Benzothiazolone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.101	186.80 ng	38628600	Phenanthrene-d10	17.131

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2(3H)-Benzothiazolone	151	C7H5NOS	000934-34-9	94
2			Thieno[3,2-b]pyridin-7-ol	151	C7H5NOS	107818-20-2	72
3			1,2-Benzisothiazol-3(2H)-one	151	C7H5NOS	002634-33-5	52
4			6-Methyl-7-oxo-4,7-dihydro-triaz...	151	C5H5N5O	057250-39-2	50
5			t-Butylhydroquinone	166	C10H14O2	001948-33-0	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-12

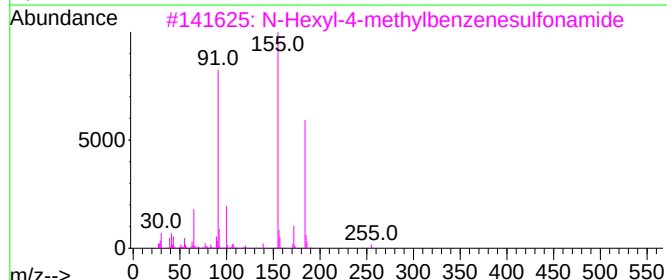
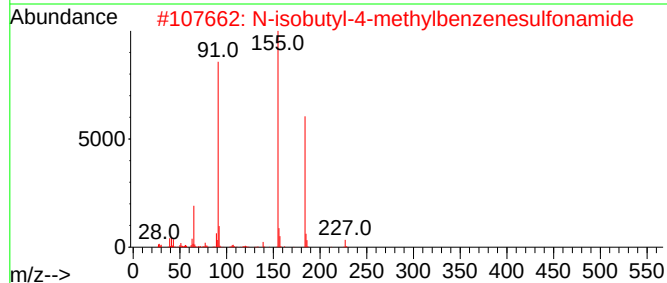
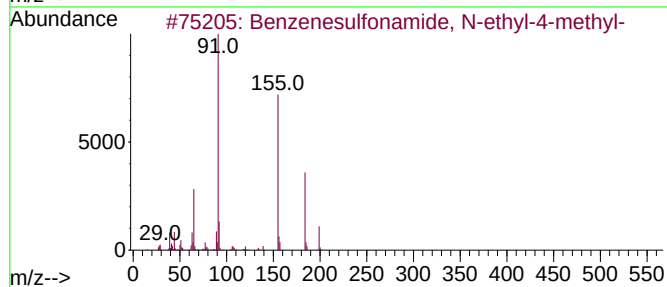
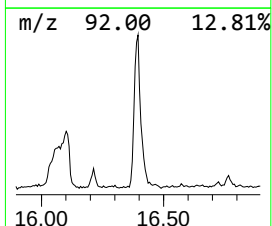
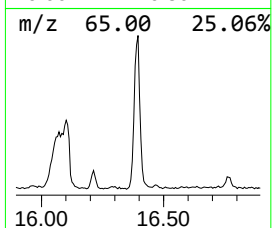
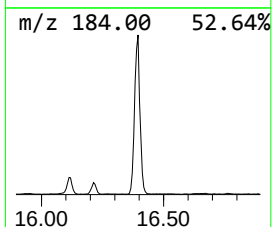
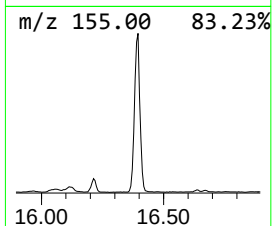
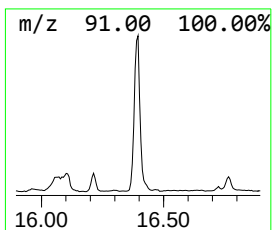
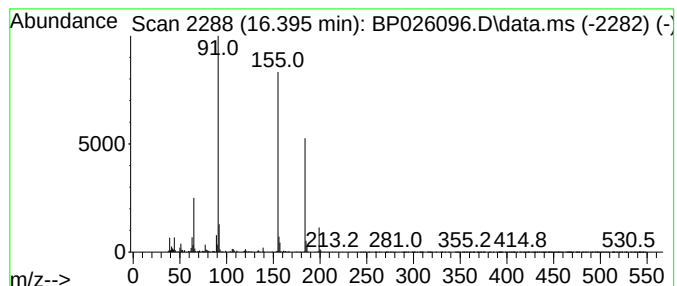
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzenesulfonamide, N-ethyl... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.395	11.28 ng	2331940	Phenanthrene-d10	17.131

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzenesulfonamide, N-ethyl-4-me...	199	C9H13NO2S	000080-39-7	97
2			N-isobutyl-4-methylbenzenesulfon...	227	C11H17NO2S	023705-38-6	87
3			N-Hexyl-4-methylbenzenesulfonamide	255	C13H21NO2S	001143-01-7	86
4			4-Methyl-N-[2-(3-nitro-[1,2,4]tr...	311	C11H13N5O4S	1000301-03-5	83
5			Furazan-3-carboxylic acid, 4-ami...	325	C12H15N5O4S	1000304-69-4	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

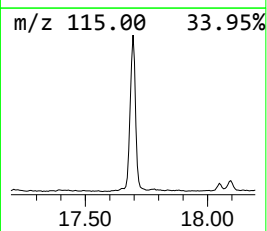
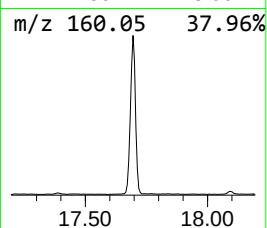
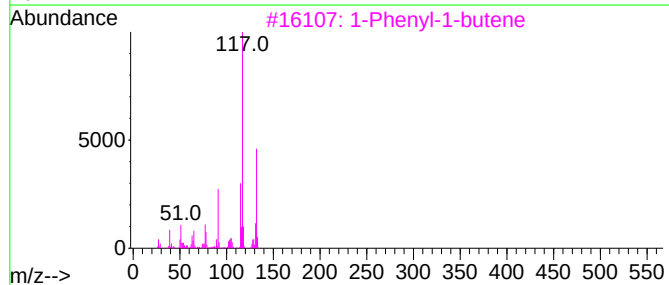
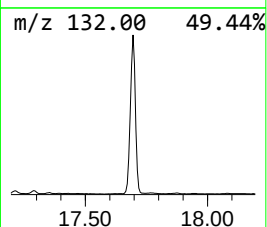
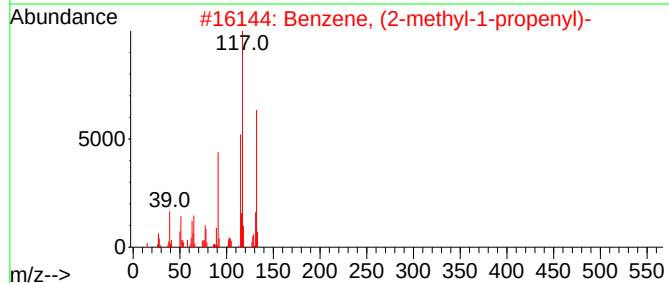
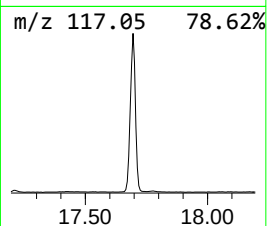
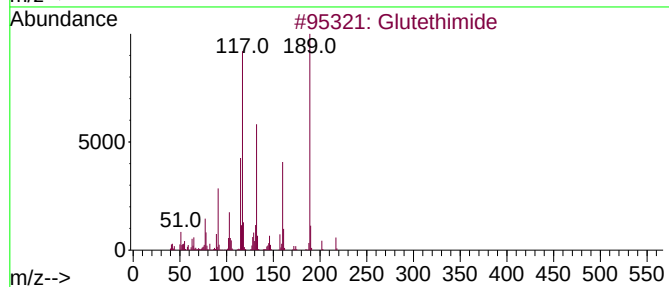
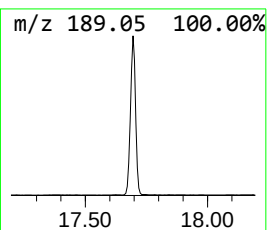
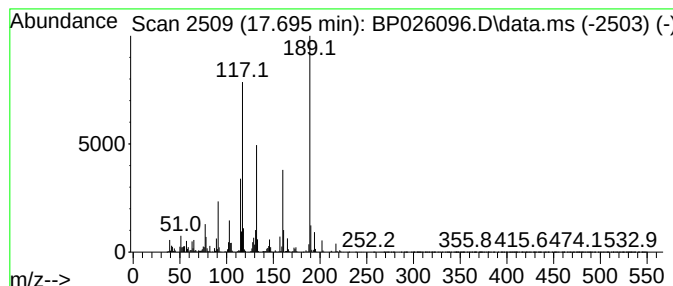
Instrument :
BNA_P
ClientSampleId :
MW-12

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 15 Glutethimide Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.695	46.49 ng	9614070	Phenanthrene-d10		17.131	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Glutethimide		217	C13H15NO2	000077-21-4	99
2	Benzene, (2-methyl-1-propenyl)-		132	C10H12	000768-49-0	89
3	1-Phenyl-1-butene		132	C10H12	000824-90-8	64
4	(E)-1-Phenyl-1-butene		132	C10H12	001005-64-7	64
5	Benzene, 2-butenyl-		132	C10H12	001560-06-1	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
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Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

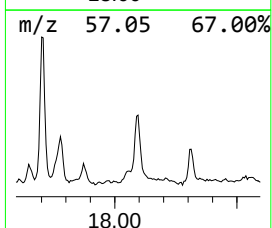
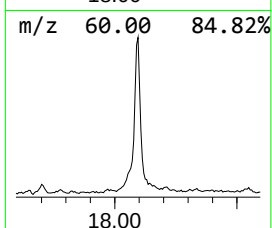
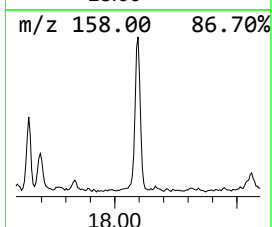
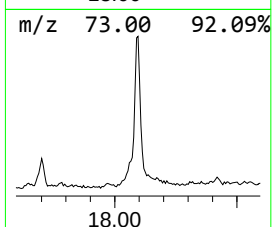
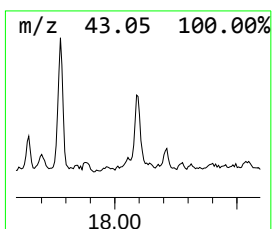
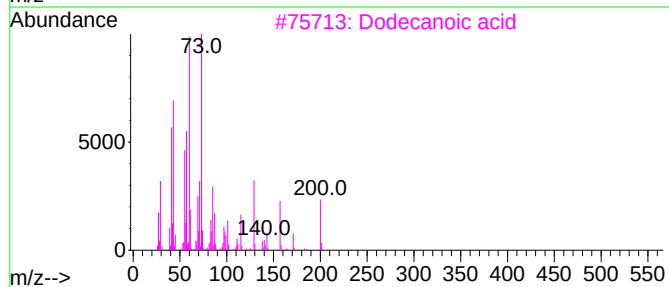
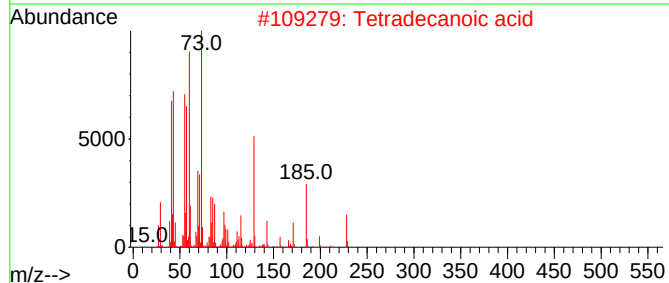
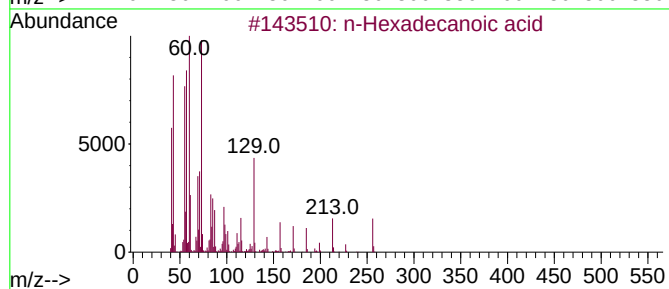
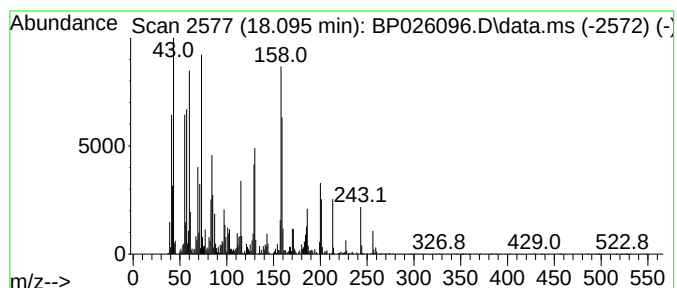
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP102925.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 17 n-Hexadecanoic acid Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.095	6.48 ng	1339110	Phenanthrene-d10	17.131	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	Tetradecanoic acid	228	C14H28O2	000544-63-8	60
3	Dodecanoic acid	200	C12H24O2	000143-07-7	50
4	1H-Indole-3-carboxaldehyde, 7-me...	159	C10H9NO	004771-50-0	38
5	1H-Indole-3-carboxaldehyde, 2-me...	159	C10H9NO	005416-80-8	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-12

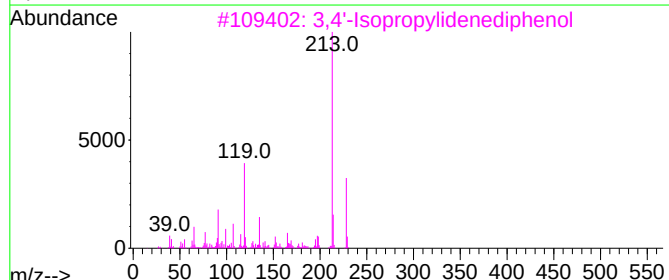
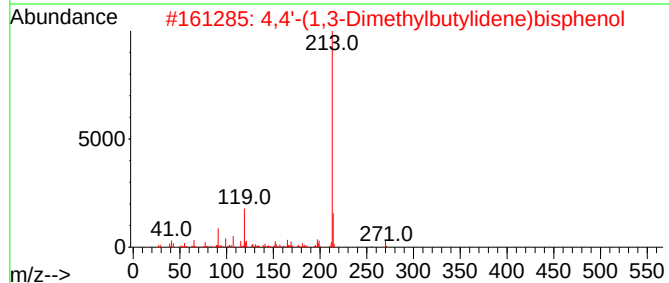
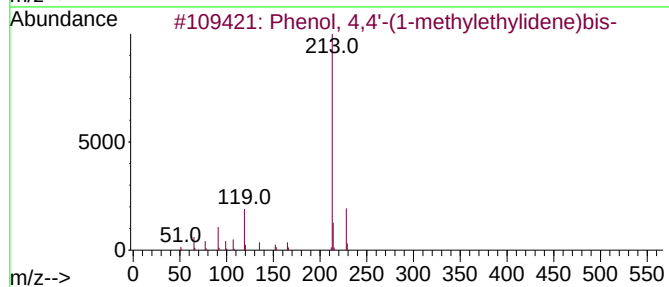
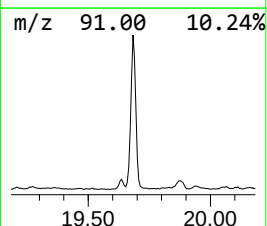
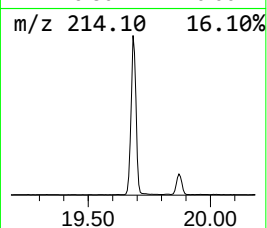
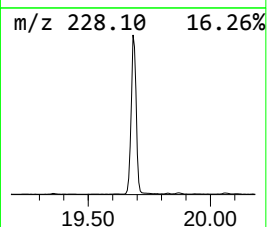
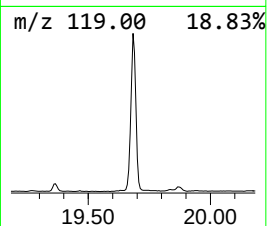
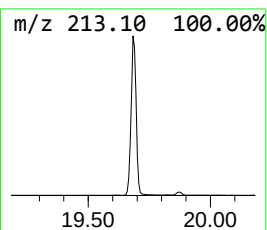
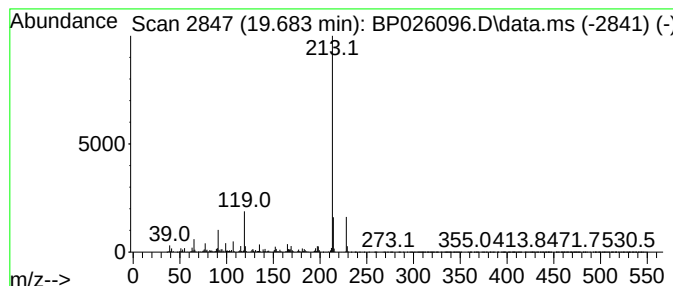
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 19 Phenol, 4,4'-(1-methylethyl... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.683	51.20 ng	13490600	Chrysene-d12	21.583

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4,4'-(1-methylethylidene...	228	C15H16O2	000080-05-7	97
2			4,4'-(1,3-Dimethylbutylidene)bis...	270	C18H22O2	006807-17-6	83
3			3,4'-Isopropylidenediphenol	228	C15H16O2	046765-25-7	70
4			2,5-bis(Trimethylsilyl)thiophene	228	C10H20SSi2	017906-71-7	59
5			5-Fluoro-2-nitrophenylamine, TMS...	228	C9H13FN2O2Si	1000484-54-5	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111125\
Data File : BP026096.D
Acq On : 11 Nov 2025 18:55
Operator : RC/JU
Sample : Q3569-02
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
MW-12

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
Bicyclo[2.2.1]h...	10.007	10.1	ng	1382570	2	10.443	2726100	20.0
Phenol, 4-chloro-	10.331	10.3	ng	1396640	2	10.443	2726100	20.0
Phenol, p-tert-...	11.796	37.4	ng	5103880	2	10.443	2726100	20.0
Benzoic acid, p...	14.142	8.7	ng	1628460	3	14.313	3764110	20.0
Ibuprofen	15.401	14.5	ng	2725660	3	14.313	3764110	20.0
2(3H)-Benzothia...	16.101	186.8	ng	38628600	4	17.131	4135760	20.0
Benzenesulfonam...	16.395	11.3	ng	2331940	4	17.131	4135760	20.0
Glutethimide	17.695	46.5	ng	9614070	4	17.131	4135760	20.0
n-Hexadecanoic ...	18.095	6.5	ng	1339110	4	17.131	4135760	20.0
Phenol, 4,4'-(1...	19.683	51.2	ng	13490600	5	21.583	5270290	20.0