

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
 Data File : BE101596.D
 Acq On : 13 Nov 2024 15:32
 Operator : RC/JU
 Sample : P4816-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 331

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_E\Methods\8270-BE110624.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BE101596.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.658	9	12	20	rVB	27696	38476	1.32%	0.188%
2	5.302	459	462	472	rVB2	31078	49875	1.71%	0.243%
3	6.935	735	740	750	rBV	51917	103388	3.54%	0.504%
4	7.546	840	844	849	rVB	39140	57061	1.95%	0.278%
5	8.757	1044	1050	1072	rVB	115145	232965	7.97%	1.136%
6	10.314	1309	1315	1322	rBV	215327	385949	13.20%	1.882%
7	11.577	1526	1530	1538	rVB3	22851	42995	1.47%	0.210%
8	12.176	1626	1632	1636	rVB6	18346	35623	1.22%	0.174%
9	12.435	1665	1676	1681	rBV6	11007	32130	1.10%	0.157%
10	12.487	1681	1685	1691	rVV6	21044	38080	1.30%	0.186%
11	12.546	1691	1695	1702	rVB3	30291	46938	1.61%	0.229%
12	12.781	1727	1735	1742	rBV	953760	1458234	49.88%	7.109%
13	12.846	1742	1746	1751	rVB3	43506	75041	2.57%	0.366%
14	13.480	1848	1854	1855	rBV6	31629	43675	1.49%	0.213%
15	13.604	1873	1875	1882	rVB6	22249	43860	1.50%	0.214%
16	13.804	1905	1909	1913	rBV5	29288	46615	1.59%	0.227%
17	13.892	1917	1924	1928	rBV	59568	115533	3.95%	0.563%
18	14.162	1962	1970	1975	rBV	463297	696810	23.84%	3.397%
19	14.221	1976	1980	1984	rVB2	35001	62480	2.14%	0.305%
20	15.531	2200	2203	2207	rVB	81598	91492	3.13%	0.446%
21	15.672	2222	2227	2233	rVB	772819	1061946	36.33%	5.177%
22	16.553	2374	2377	2383	rVB2	139719	192526	6.59%	0.939%
23	16.900	2432	2436	2440	rBV	507022	726909	24.87%	3.544%
24	16.941	2440	2443	2454	rVB	409550	639508	21.88%	3.118%
25	18.463	2699	2702	2705	rBV	175671	218369	7.47%	1.065%
26	18.668	2734	2737	2741	rVB	169314	204679	7.00%	0.998%
27	18.956	2781	2786	2794	rVB	1136991	1568934	53.67%	7.649%
28	19.321	2844	2848	2853	rVB	1001967	1323553	45.28%	6.452%
29	19.497	2873	2878	2891	rVB	2102185	2923337	100.00%	14.252%
30	20.149	2986	2989	2995	rVB2	200466	286656	9.81%	1.397%
31	20.713	3082	3085	3089	rBV2	176337	221280	7.57%	1.079%
32	21.072	3139	3146	3149	rBV2	771619	1729795	59.17%	8.433%
33	21.101	3149	3151	3156	rVB	638780	813787	27.84%	3.967%
34	22.323	3355	3359	3363	rVB	252912	362048	12.38%	1.765%
35	22.646	3410	3414	3428	rVB	545664	1841454	62.99%	8.977%
36	23.151	3496	3500	3509	rVB	392318	844269	28.88%	4.116%

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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	23.245	3512	3516	3527	rVB	397663	800257	27.37%	3.901%
38	23.357	3530	3535	3540	rBV	530032	1055887	36.12%	5.148%

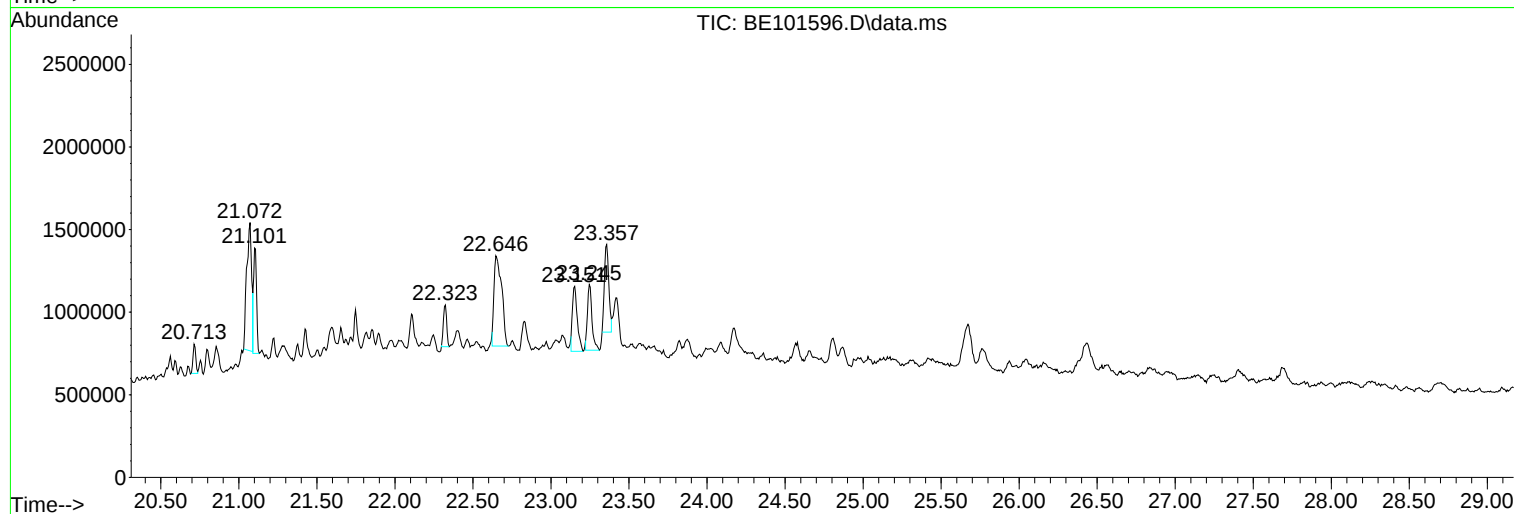
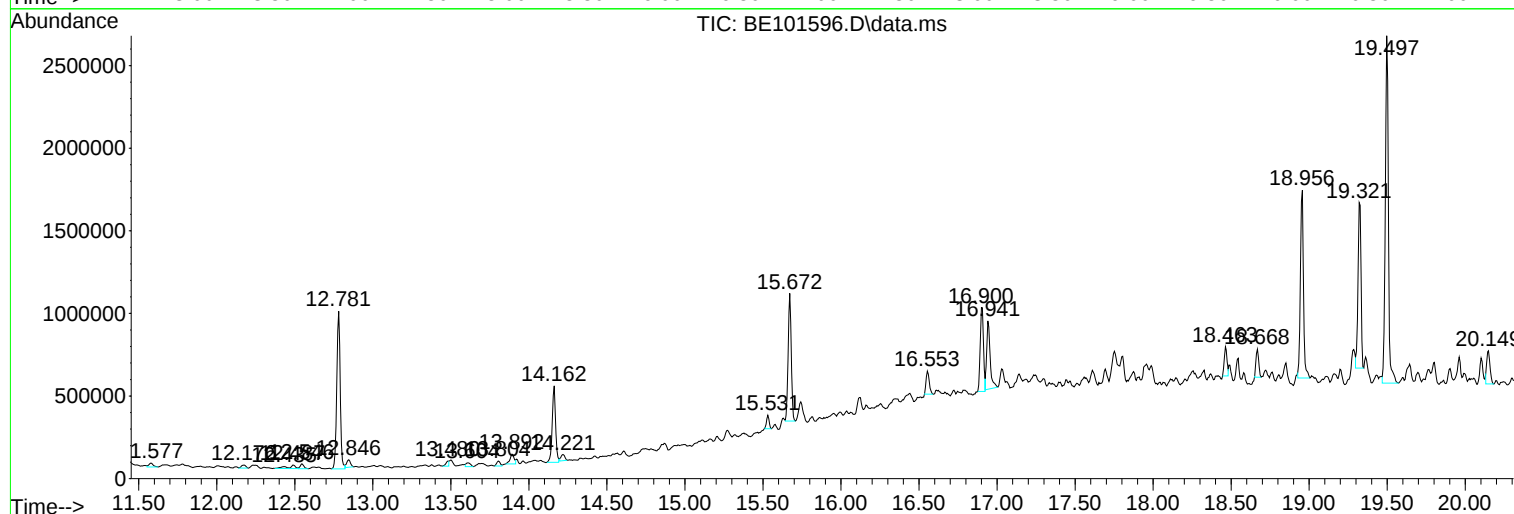
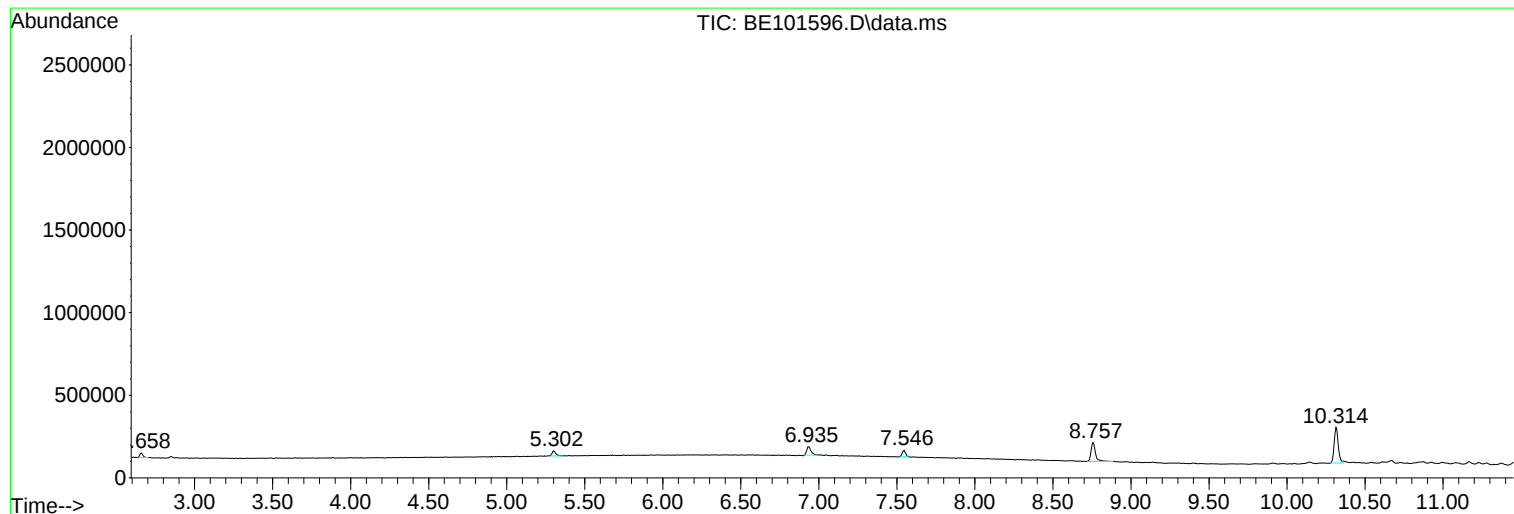
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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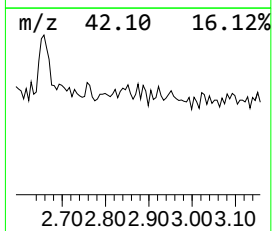
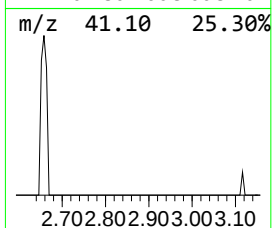
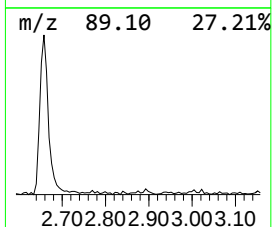
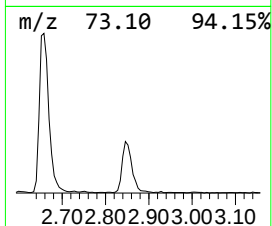
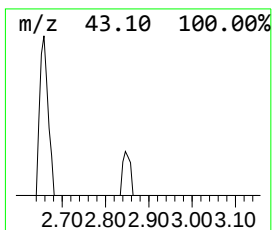
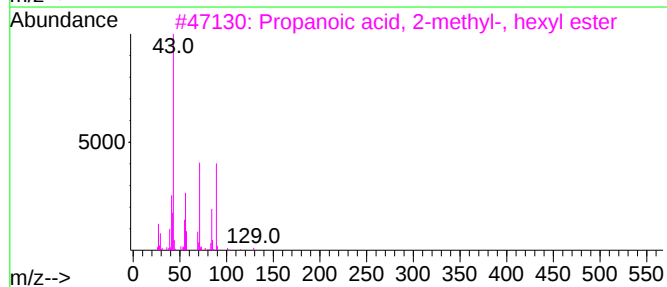
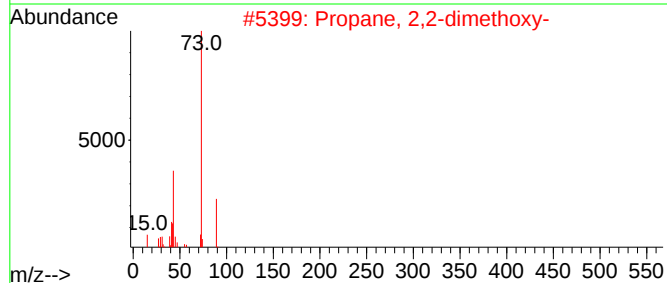
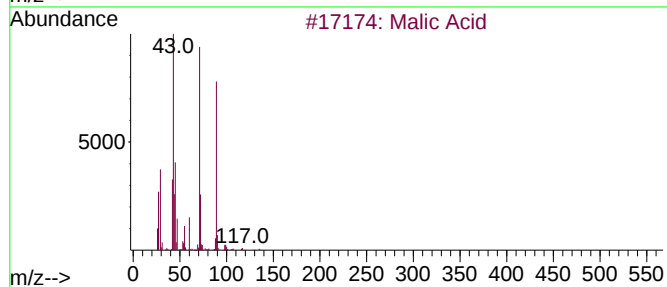
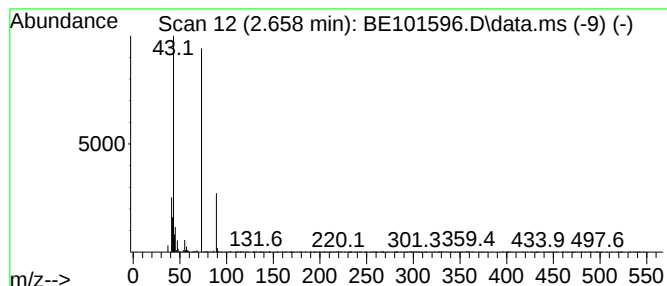
Instrument :
BNA_E
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 unknown2.658 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
2.658	13.49 ng	38476	1,4-Dichlorobenzene-d4	7.546	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Malic Acid	134	C4H6O5	006915-15-7	27
2	Propane, 2,2-dimethoxy-	104	C5H12O2	000077-76-9	17
3	Propanoic acid, 2-methyl-, hexyl...	172	C10H20O2	002349-07-7	12
4	2-Propanone, 1,1,3,3-tetrapropoxy-	290	C15H30O5	113358-60-4	12
5	1,3-Dioxolane, 2-heptyl-	172	C10H20O2	004359-57-3	9



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Misc :
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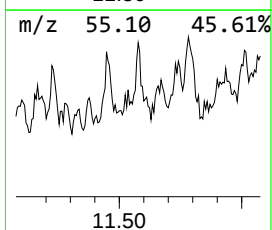
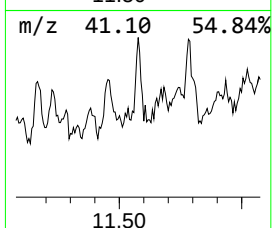
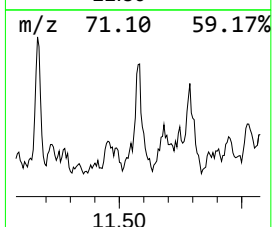
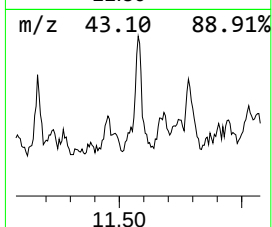
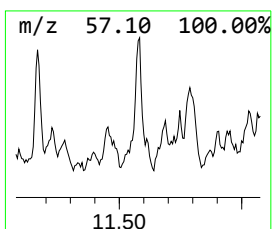
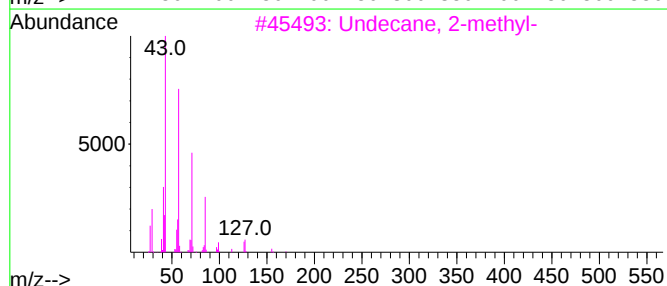
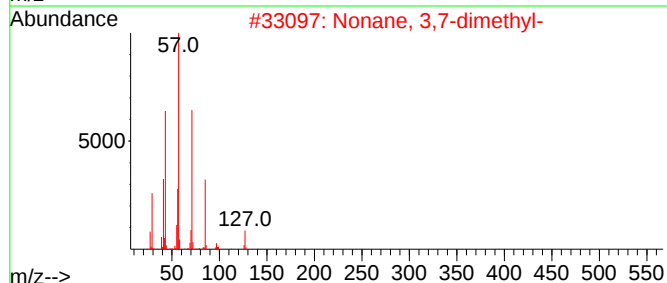
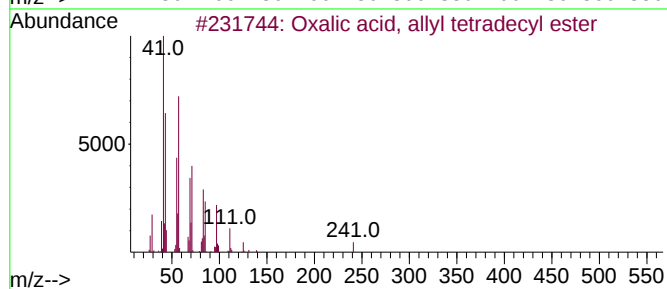
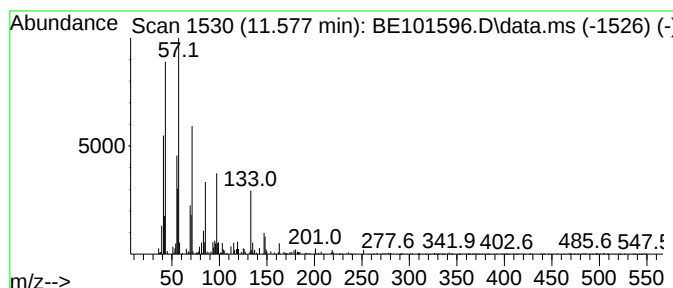
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Oxalic acid, allyl tetradec... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.577	2.23 ng	42995	Naphthalene-d8	10.313	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Oxalic acid, allyl tetradecyl ester	326	C19H34O4	1000309-24-2	52
2	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	43
3	Undecane, 2-methyl-	170	C12H26	007045-71-8	43
4	Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	38
5	Decane, 2,3,5-trimethyl-	184	C13H28	062238-11-3	38



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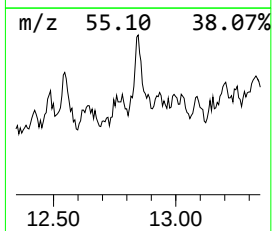
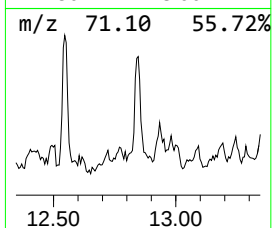
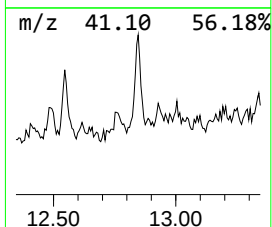
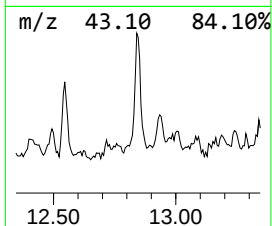
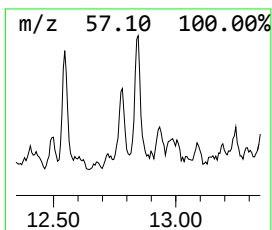
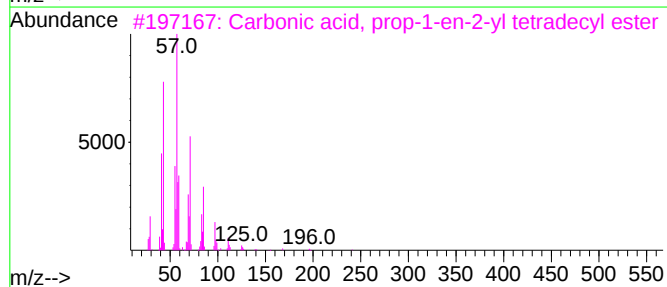
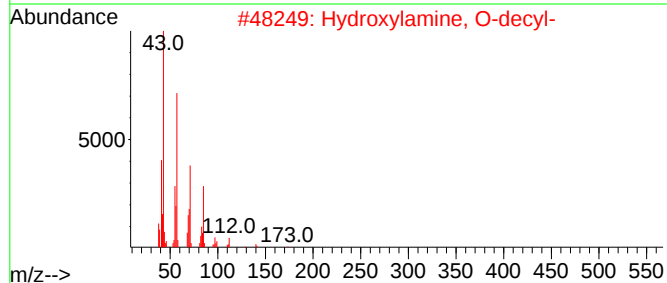
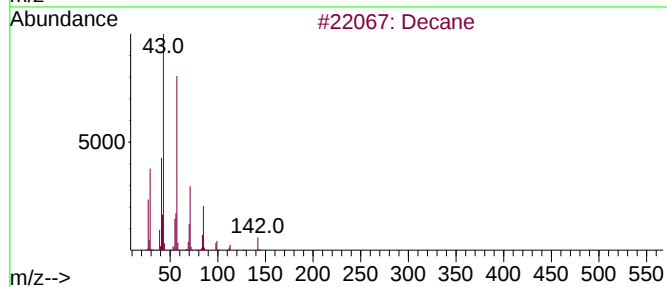
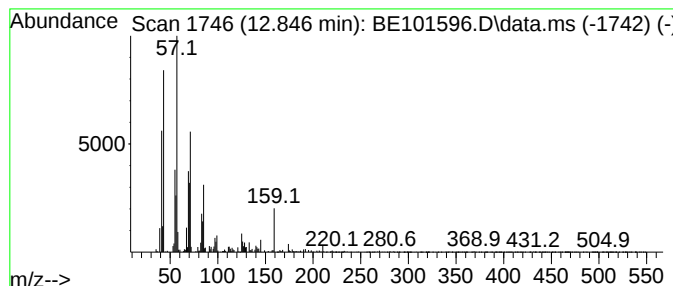
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Decane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
12.846	2.15 ng	75041	Acenaphthene-d10	14.162	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	76
2	Hydroxylamine, O-decyl-	173	C10H23NO	029812-79-1	72
3	Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	64
4	Tetradecane	198	C14H30	000629-59-4	52
5	Heneicosane	296	C21H44	000629-94-7	52



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Misc :
ALS Vial : 5 Sample Multiplier: 1

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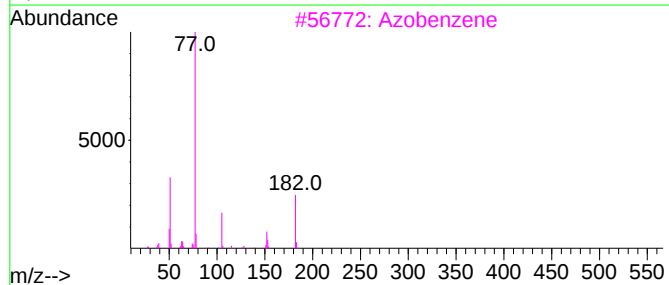
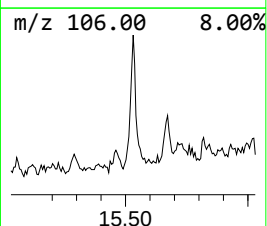
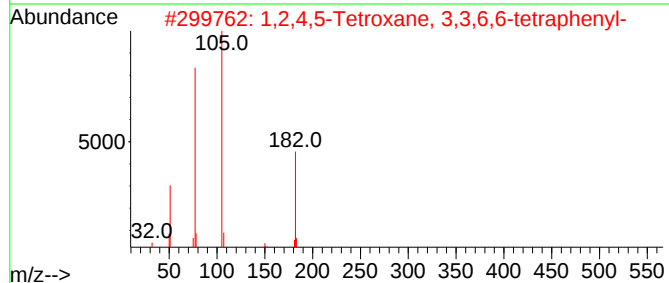
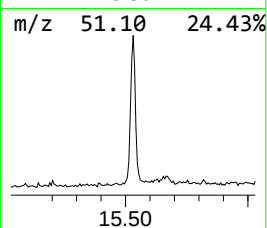
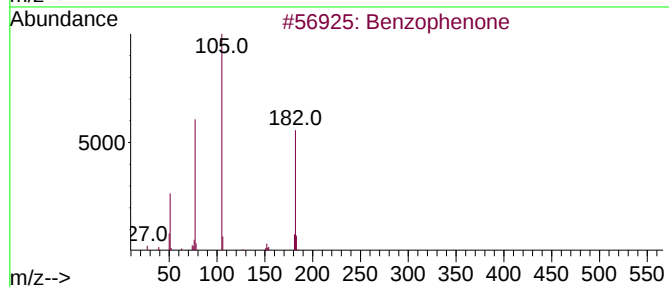
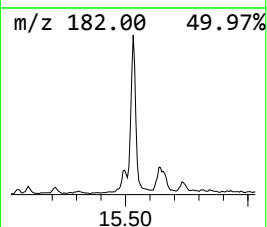
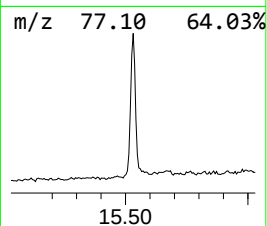
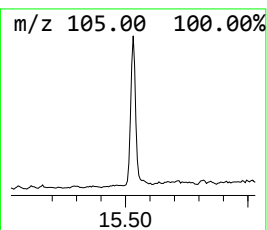
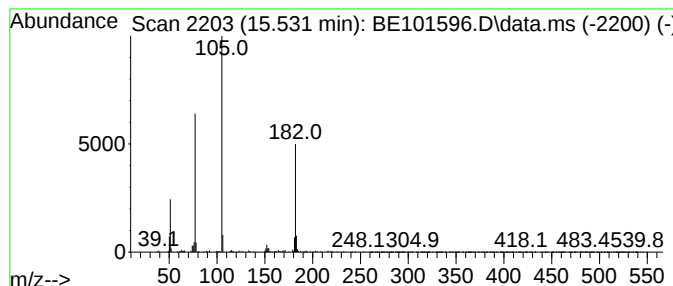
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzophenone Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.531	2.63 ng	91492	Acenaphthene-d10	14.162

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64
3			Azobenzene	182	C12H10N2	000103-33-3	59
4			Benzoic acid, 2-benzoyl-	226	C14H10O3	000085-52-9	56
5			Methanone, phenyl-2-pyridinyl-	183	C12H9NO	000091-02-1	56



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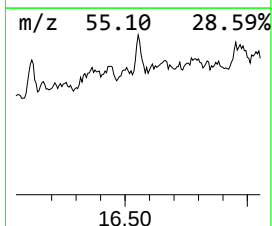
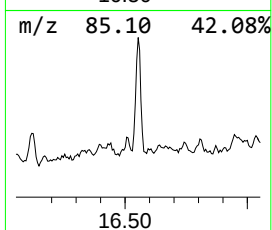
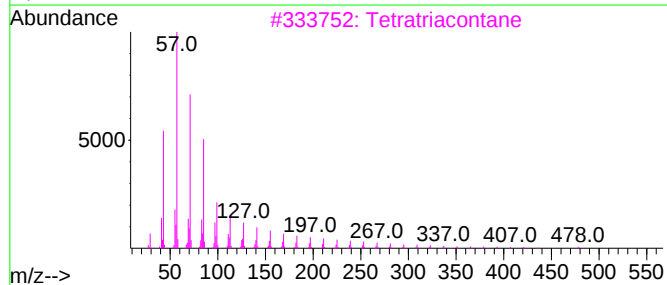
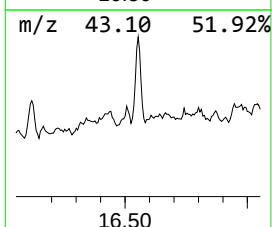
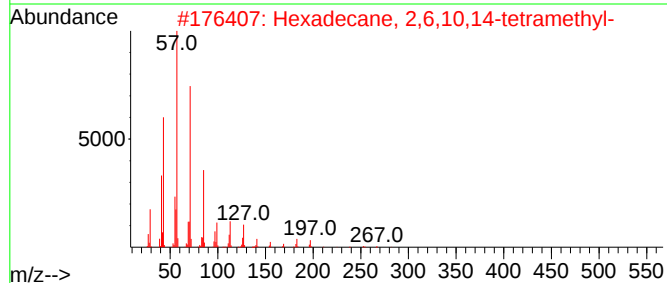
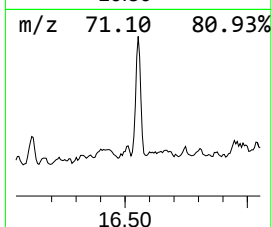
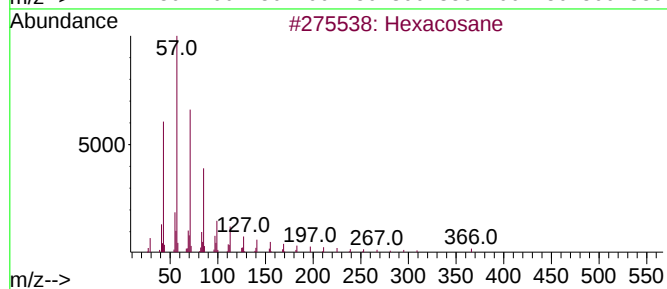
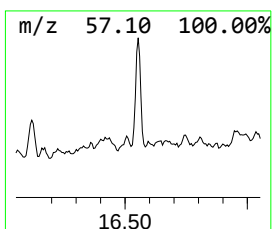
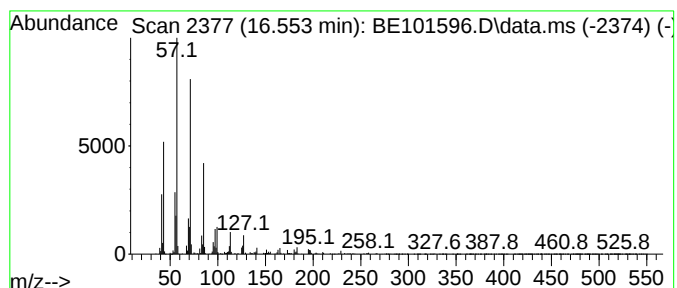
Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.553	5.30 ng	192526	Phenanthrene-d10	16.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	90
2			Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
3			Tetratriacontane	479	C34H70	014167-59-0	86
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	86
5			Disulfide, di-tert-dodecyl	402	C24H50S2	027458-90-8	83



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Data File : BE101596.D
Acq On : 13 Nov 2024 15:32
Operator : RC/JU
Sample : P4816-01
Misc :
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Instrument :
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ClientSampleId :
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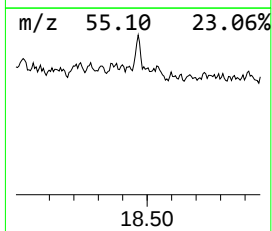
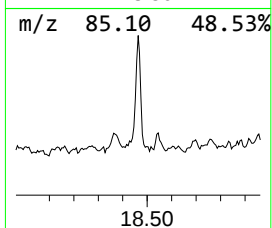
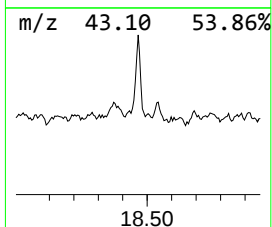
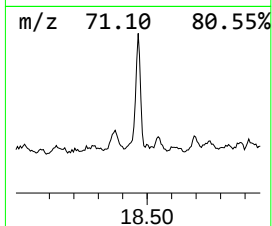
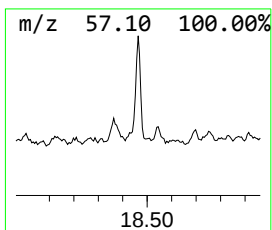
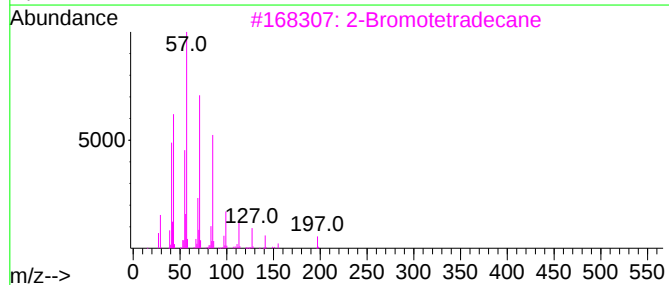
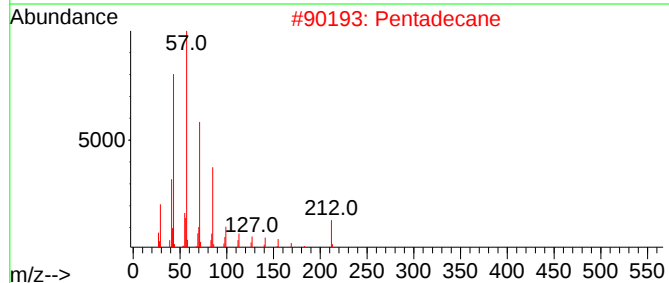
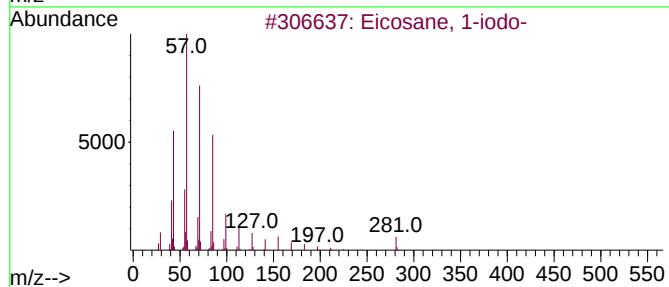
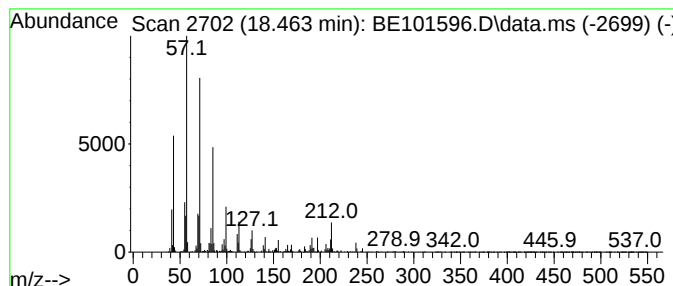
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Eicosane, 1-iodo- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.463	6.01 ng	218369	Phenanthrene-d10	16.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 1-iodo-	408	C20H41I	1000406-31-8	87
2			Pentadecane	212	C15H32	000629-62-9	83
3			2-Bromotetradecane	276	C14H29Br	074036-95-6	83
4			Octadecane, 1-iodo-	380	C18H37I	000629-93-6	83
5			3-Ethyl-2,6,10-trimethylundecane	226	C16H34	1000432-25-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
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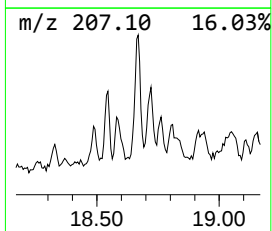
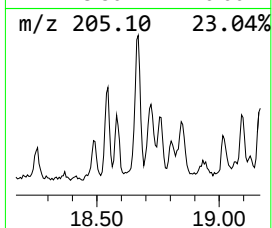
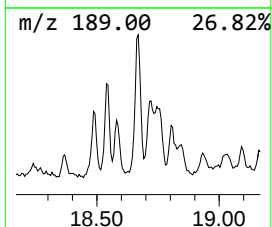
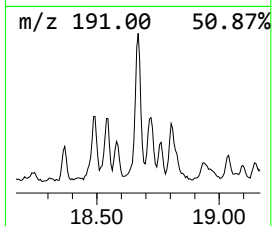
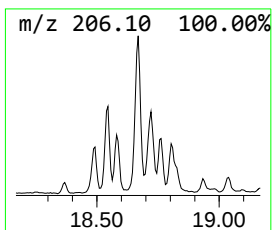
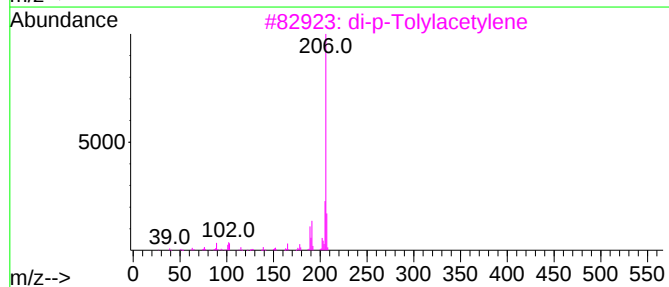
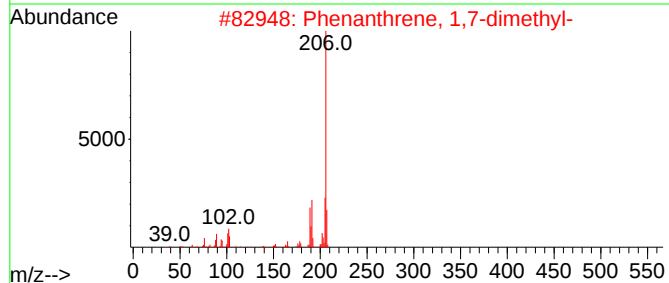
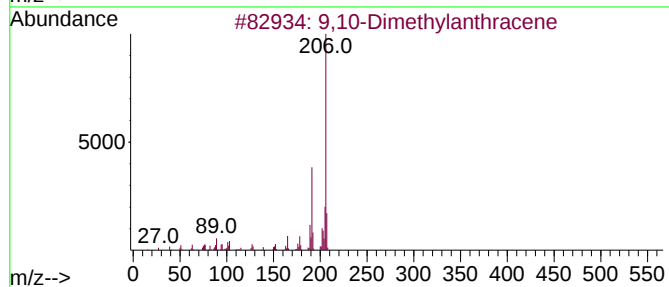
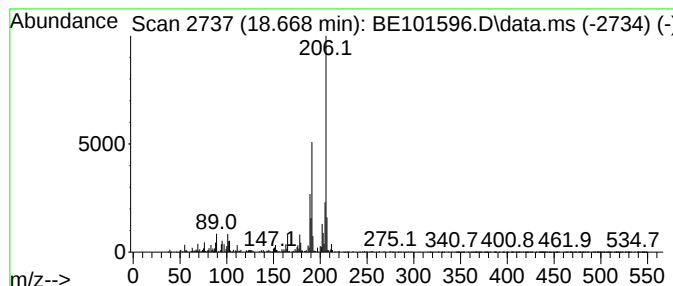
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 9,10-Dimethylantracene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.669	5.63 ng	204679	Phenanthrene-d10	16.906

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9,10-Dimethylantracene	206	C16H14	000781-43-1	94
2			Phenanthrene, 1,7-dimethyl-	206	C16H14	000483-87-4	93
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	91
4			1H-Indene, 2-methyl-3-phenyl-	206	C16H14	035099-60-6	90
5			3-Methyl-1-phenyl-1H-indene	206	C16H14	022360-63-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
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Sample : P4816-01
Misc :
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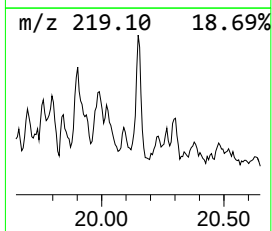
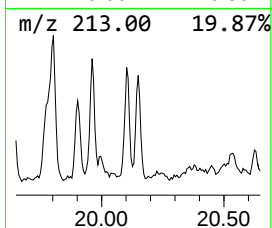
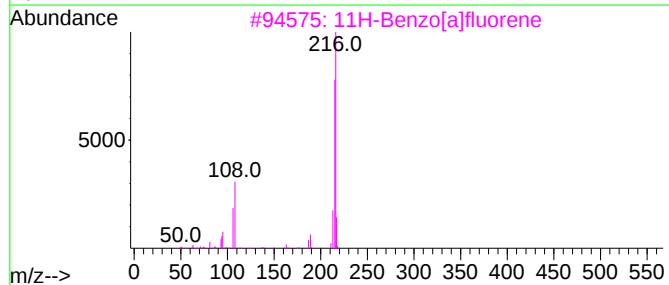
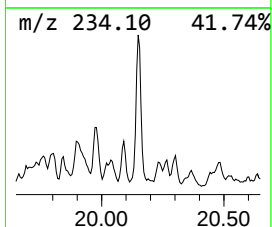
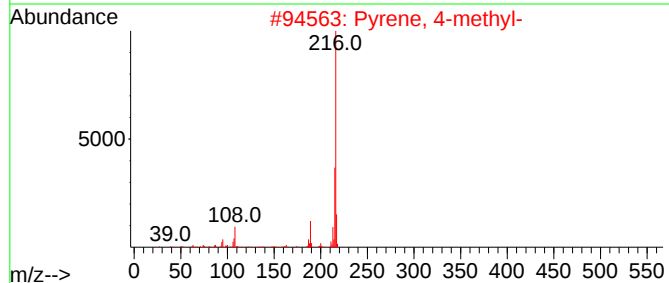
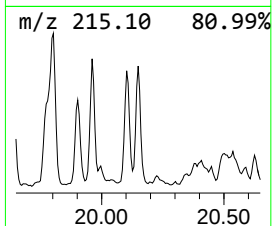
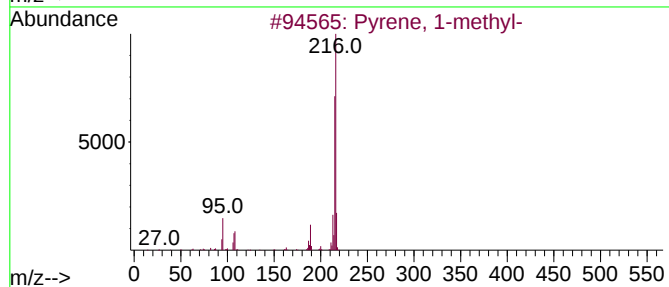
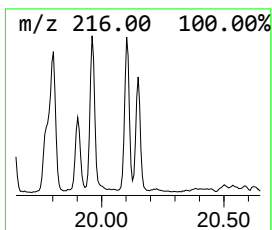
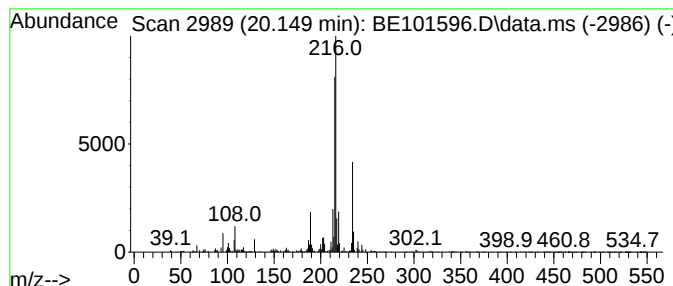
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Pyrene, 1-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.149	3.31 ng	286656	Chrysene-d12	21.072

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	97
2			Pyrene, 4-methyl-	216	C17H12	003353-12-6	80
3			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	62
4			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	60
5			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
Acq On : 13 Nov 2024 15:32
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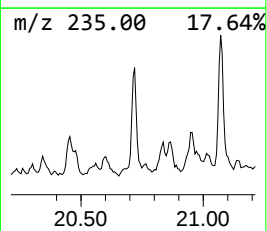
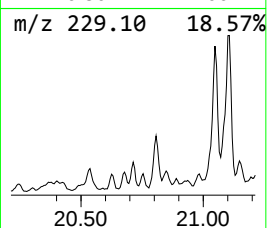
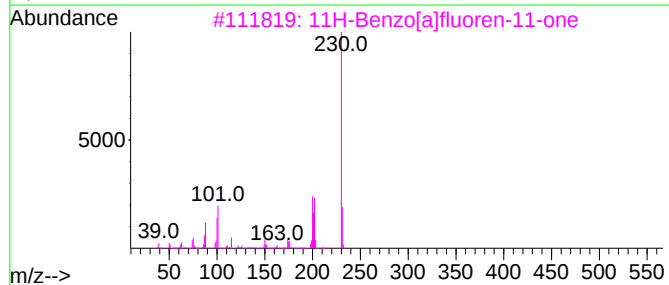
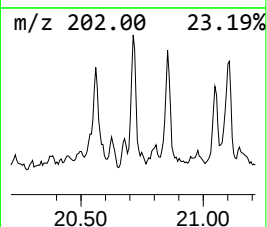
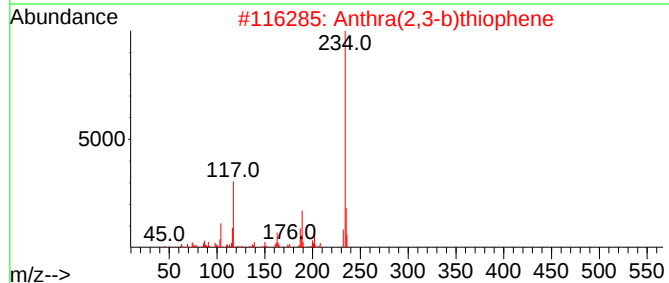
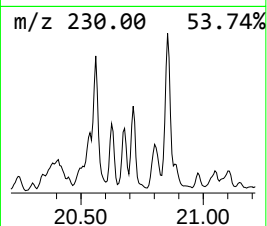
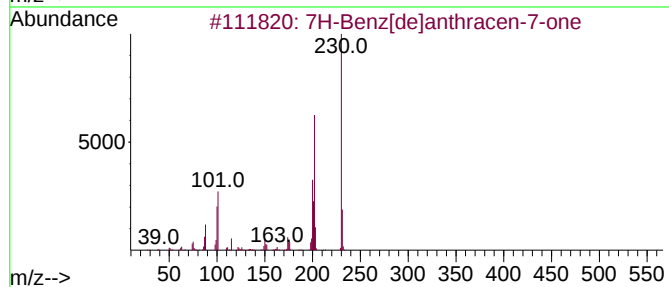
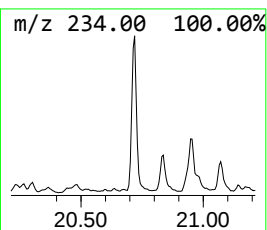
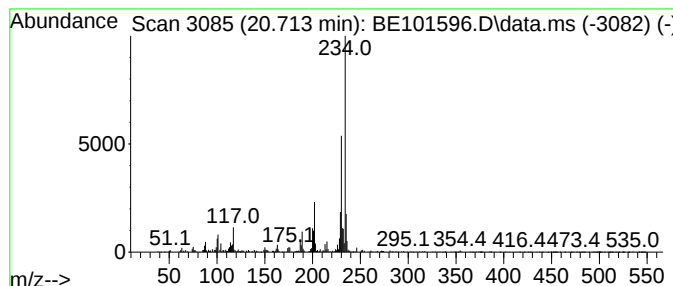
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 7H-Benz[de]anthracen-7-one Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.713	2.56 ng	221280	Chrysene-d12	21.072

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	93
2			Anthra(2,3-b)thiophene	234	C16H10S	022108-55-0	89
3			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	74
4			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	60
5			N-(4-Methylphenyl)quinolin-4-amine	234	C16H14N2	1000484-32-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
Acq On : 13 Nov 2024 15:32
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Sample : P4816-01
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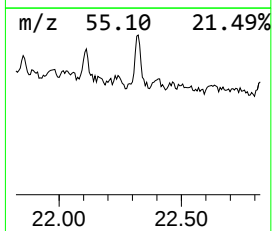
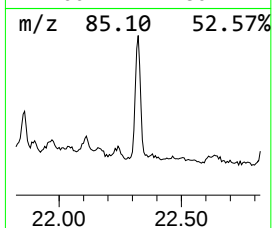
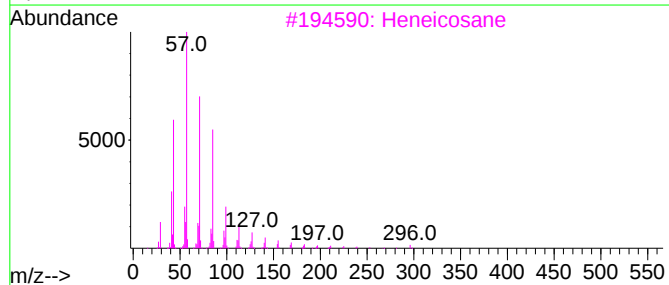
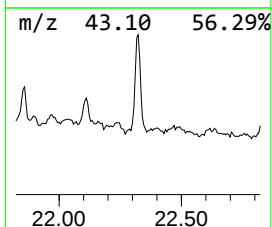
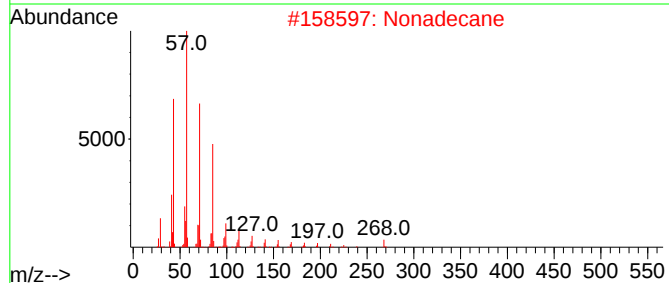
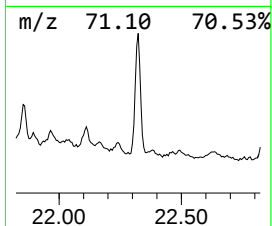
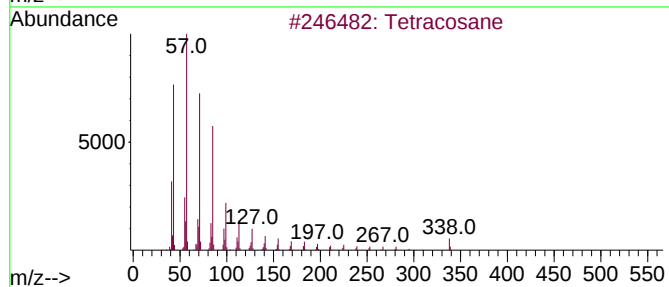
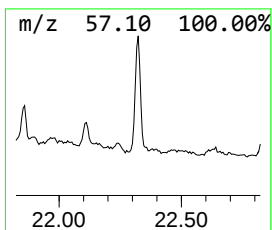
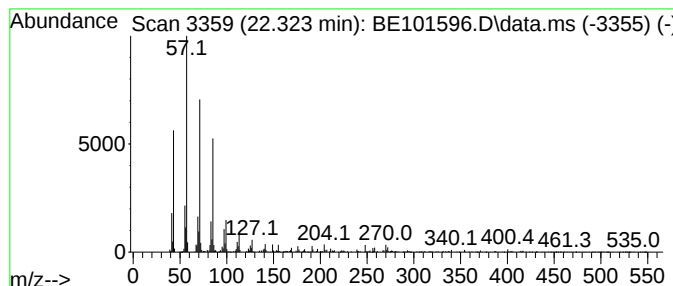
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tetracosane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.323	6.86 ng	362048	Perylene-d12	23.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	96
2			Nonadecane	268	C19H40	000629-92-5	93
3			Heneicosane	296	C21H44	000629-94-7	90
4			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	90
5			Dotriacontane	451	C32H66	000544-85-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
Acq On : 13 Nov 2024 15:32
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ALS Vial : 5 Sample Multiplier: 1

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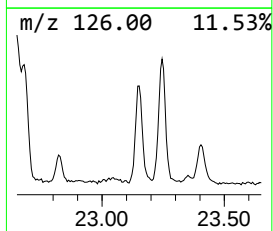
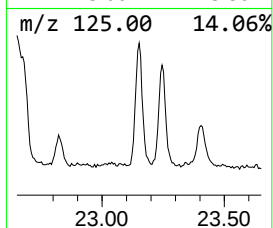
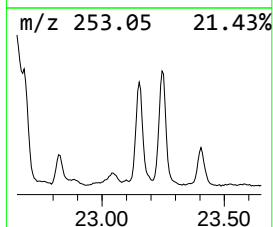
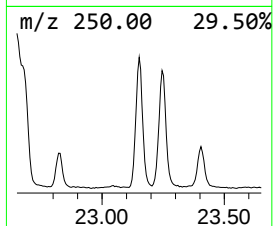
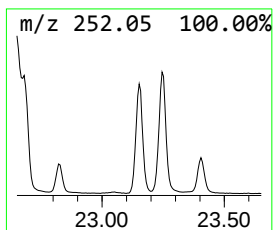
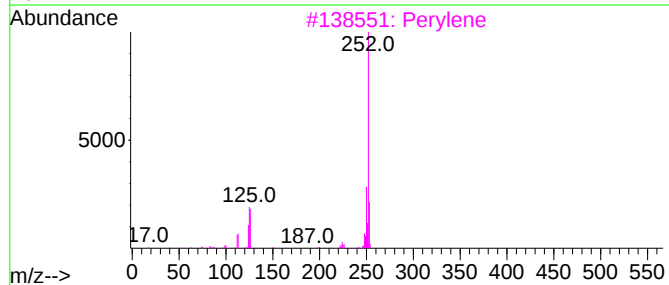
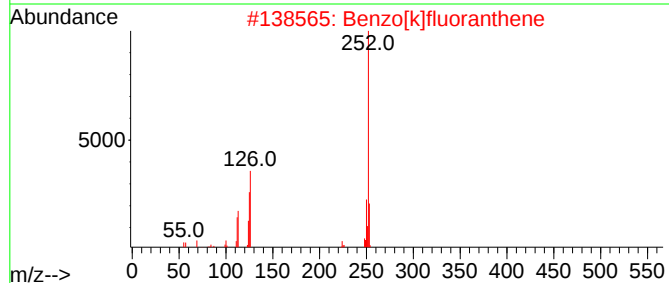
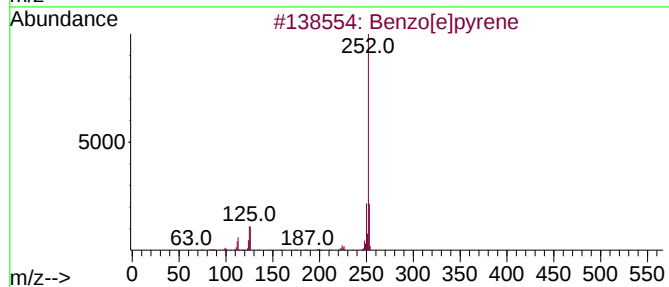
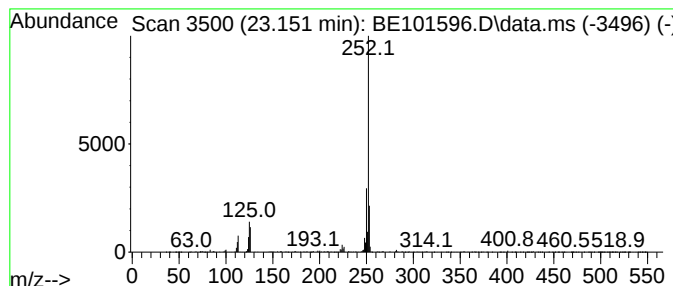
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.151	15.99 ng	844269	Perylene-d12	23.357

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
3			Perylene	252	C20H12	000198-55-0	96
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	94
5			Benzo[a]pyrene	252	C20H12	000050-32-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE111324\
Data File : BE101596.D
Acq On : 13 Nov 2024 15:32
Operator : RC/JU
Sample : P4816-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_E\Methods\8270-BE110624.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
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Decane	12.846	2.1	ng	75041	3	14.162	696810	20.0
Benzophenone	15.531	2.6	ng	91492	3	14.162	696810	20.0
Hexacosane	16.553	5.3	ng	192526	4	16.906	726909	20.0
Eicosane, 1-iodo-	18.463	6.0	ng	218369	4	16.906	726909	20.0
9,10-Dimethylan...	18.669	5.6	ng	204679	4	16.906	726909	20.0
Pyrene, 1-methyl-	20.149	3.3	ng	286656	5	21.072	1729800	20.0
7H-Benz[de]anth...	20.713	2.6	ng	221280	5	21.072	1729800	20.0
Tetracosane	22.323	6.9	ng	362048	6	23.357	1055890	20.0
Benzo[e]pyrene	23.151	16.0	ng	844269	6	23.357	1055890	20.0