

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
 Data File : BF144117.D
 Acq On : 28 Oct 2025 23:10
 Operator : RC/JU
 Sample : Q3395-05
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PR132-S07-102116-20251017

Integration Parameters: rteint.p

Integrator: RTE

Smoothing : ON

Filtering: 5

Sampling : 1

Min Area: 3 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF144117.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.122	493	498	507	rVB	19287	34653	1.66%	0.197%
2	5.504	556	563	574	rBV	1361255	1894502	90.69%	10.761%
3	6.504	727	733	748	rBV	1367178	2016032	96.50%	11.452%
4	6.875	791	796	803	rVB	320064	413031	19.77%	2.346%
5	7.439	881	892	901	rBV	919432	1257984	60.22%	7.146%
6	7.692	931	935	943	rVB2	16029	26018	1.25%	0.148%
7	8.157	1009	1014	1021	rBV	386447	526517	25.20%	2.991%
8	8.216	1021	1024	1027	rVB2	18361	21260	1.02%	0.121%
9	8.375	1046	1051	1059	rBV3	15897	33448	1.60%	0.190%
10	8.575	1081	1085	1096	rVB	31987	62899	3.01%	0.357%
11	9.175	1180	1187	1192	rBV	38986	58738	2.81%	0.334%
12	9.233	1192	1197	1203	rBV	1583882	2089073	100.00%	11.867%
13	9.316	1207	1211	1215	rBV6	17917	26604	1.27%	0.151%
14	9.633	1260	1265	1272	rBV3	56006	91359	4.37%	0.519%
15	9.780	1286	1290	1294	rBV3	15850	23475	1.12%	0.133%
16	9.822	1294	1297	1303	rVB5	18285	27032	1.29%	0.154%
17	9.916	1308	1313	1318	rVV	467427	578179	27.68%	3.284%
18	10.169	1352	1356	1364	rVB6	18550	32728	1.57%	0.186%
19	10.322	1379	1382	1388	rBV6	15102	23250	1.11%	0.132%
20	10.563	1418	1423	1428	rBV	49082	73245	3.51%	0.416%
21	10.633	1429	1435	1441	rBV	283290	370317	17.73%	2.104%
22	10.710	1442	1448	1458	rVV	1142836	1483590	71.02%	8.427%
23	10.827	1464	1468	1475	rVB	82437	117125	5.61%	0.665%
24	10.939	1483	1487	1495	rVB	60031	75332	3.61%	0.428%
25	11.298	1543	1548	1554	rVB	47559	75375	3.61%	0.428%
26	11.410	1562	1567	1574	rBV	397070	559493	26.78%	3.178%
27	11.645	1604	1607	1611	rBV2	16799	22722	1.09%	0.129%
28	11.933	1651	1656	1663	rBV	118102	189724	9.08%	1.078%
29	12.992	1831	1836	1845	rBV	1170751	1546977	74.05%	8.787%
30	14.051	2012	2016	2032	rVB	315615	491679	23.54%	2.793%
31	15.168	2202	2206	2212	rVB	70698	103663	4.96%	0.589%
32	15.539	2264	2269	2278	rVB2	320585	653749	31.29%	3.714%
33	15.998	2342	2347	2354	rBV	183319	330190	15.81%	1.876%
34	16.515	2429	2435	2444	rVB	228706	462449	22.14%	2.627%
35	17.127	2531	2539	2550	rBV	232372	600053	28.72%	3.408%
36	17.845	2653	2661	2672	rVB	203397	569459	27.26%	3.235%

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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_F\Methods\8270-BF102425.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	18.697	2797	2806	2829	rVB2	174475	642687	30.76%	3.651%
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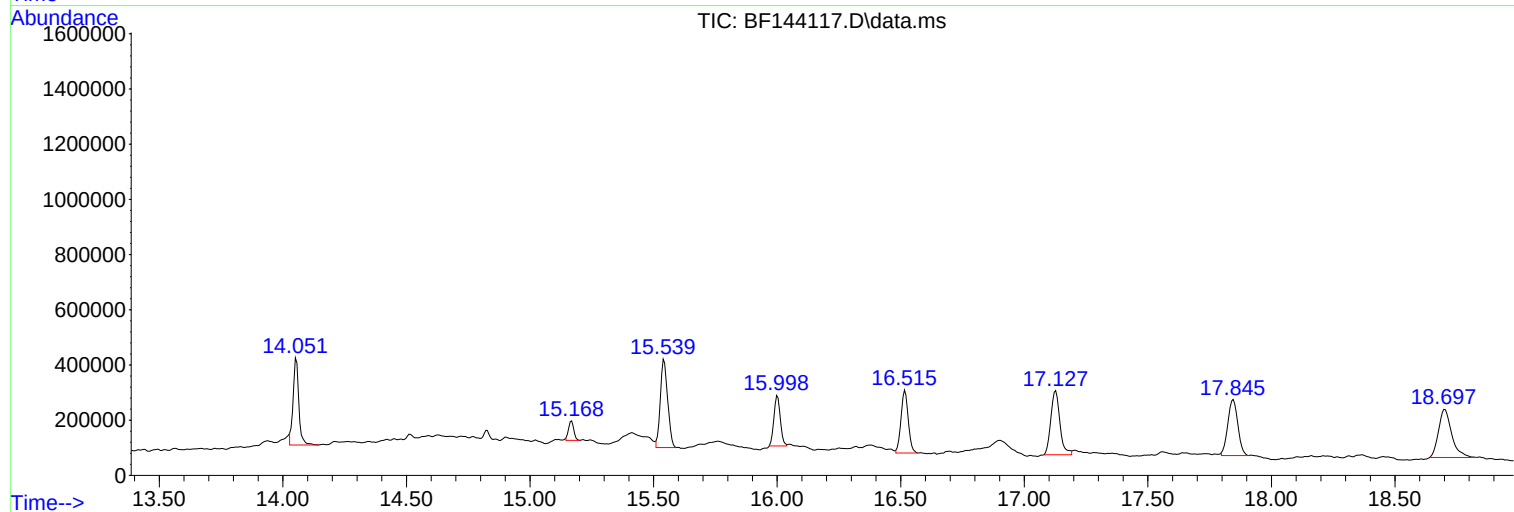
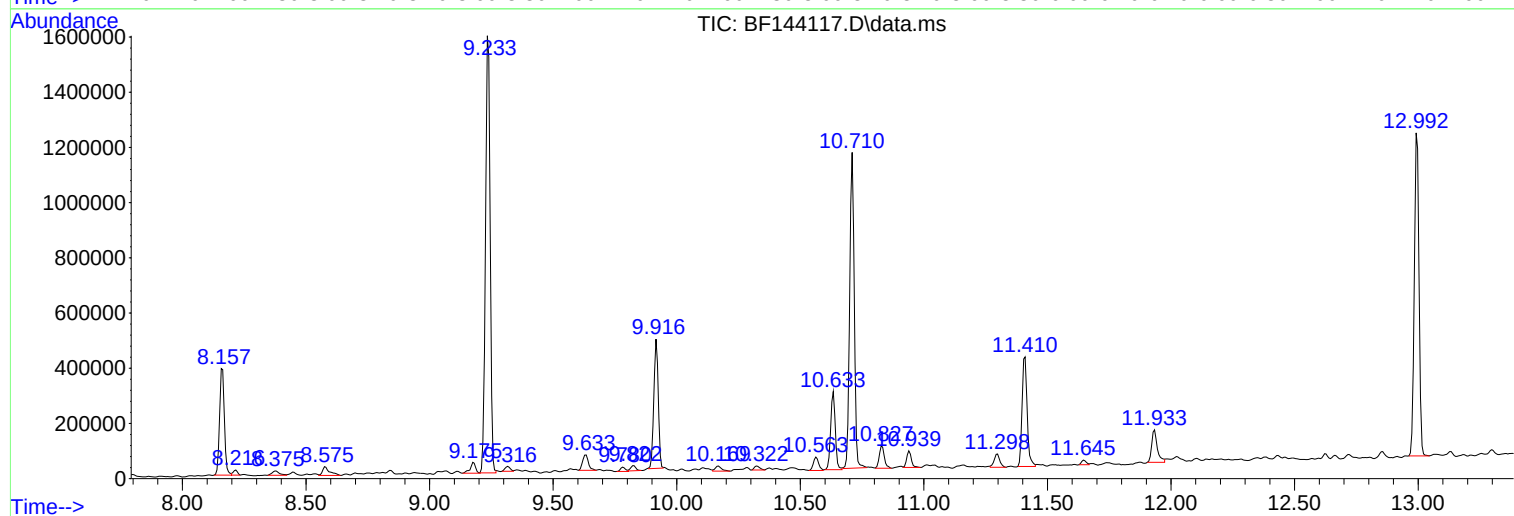
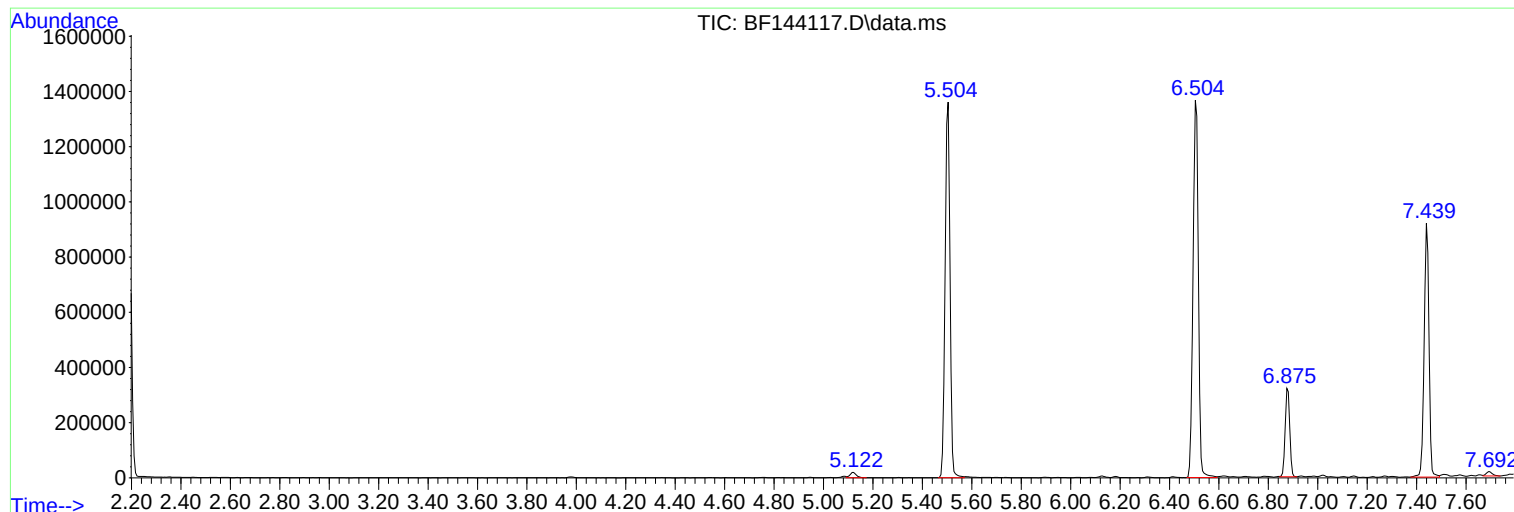
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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PR132-S07-102116-20251017

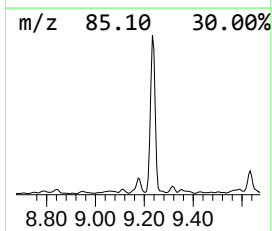
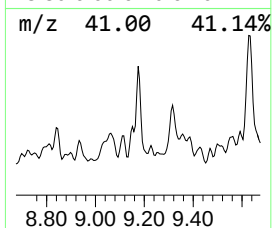
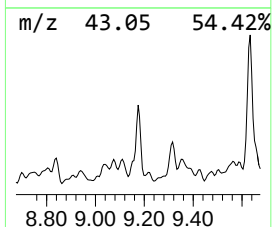
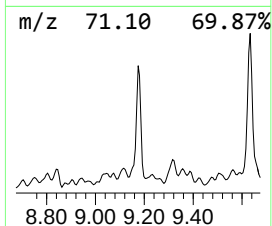
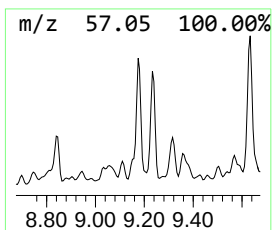
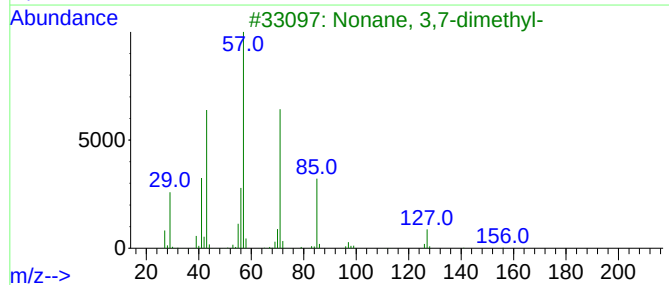
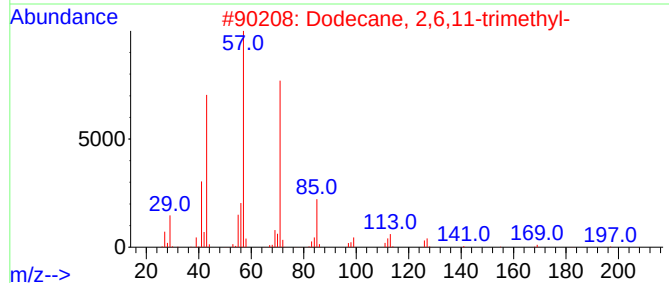
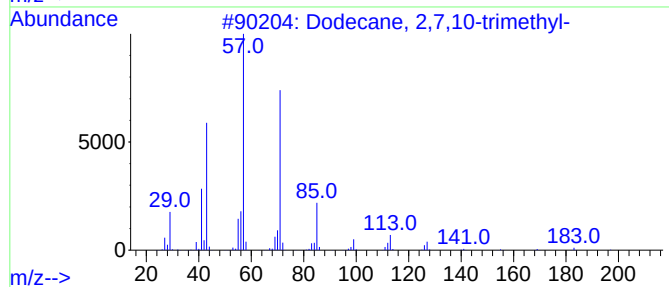
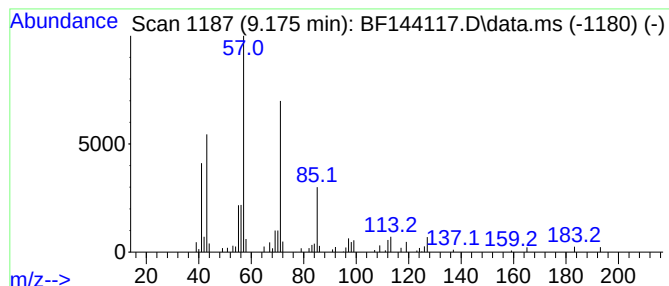
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Dodecane, 2,7,10-trimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.175	2.03 ng	58738	Acenaphthene-d10	9.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	90
2			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	72
3			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	64
4			Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	59
5			Tetracosane	338	C24H50	000646-31-1	59



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Misc :
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Instrument :
BNA_F
ClientSampleId :
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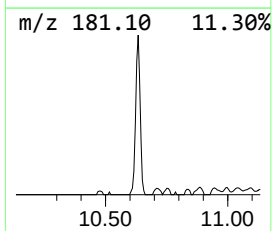
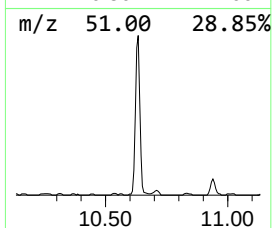
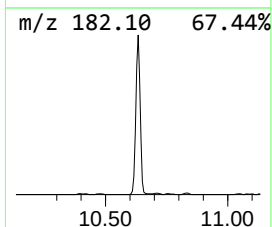
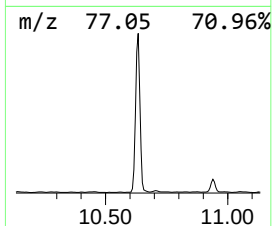
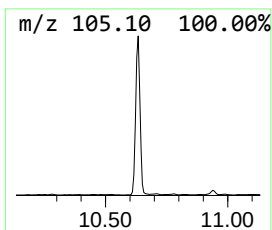
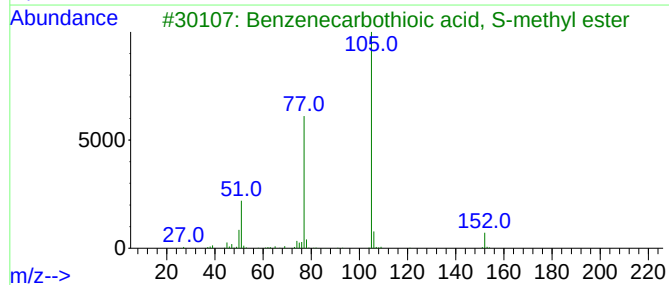
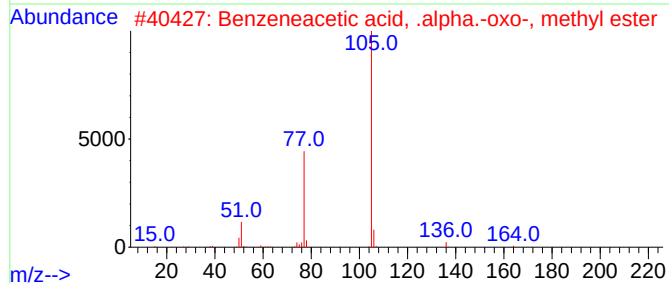
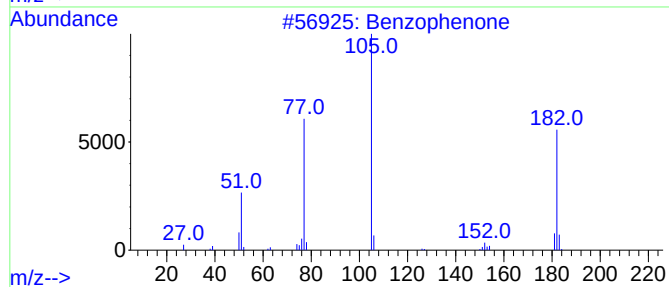
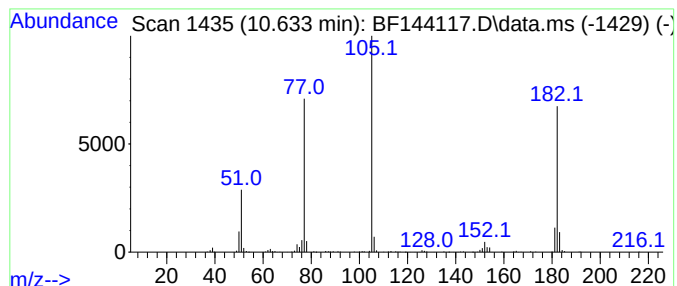
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TIC Library : C:\Database\NIST20.L
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Peak Number 5 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.633	12.81 ng	370317	Acenaphthene-d10	9.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzeneacetic acid, .alpha.-oxo-...	164	C9H8O3	015206-55-0	60
3			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	53
4			Benzoylformic acid	150	C8H6O3	000611-73-4	53
5			Pentafluoroethyl phenyl ketone	224	C9H5F5O	000394-52-5	47



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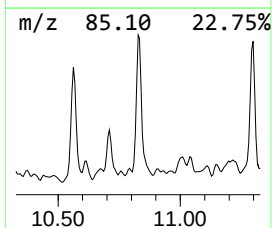
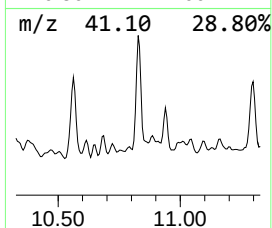
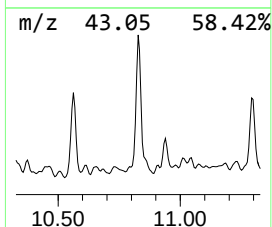
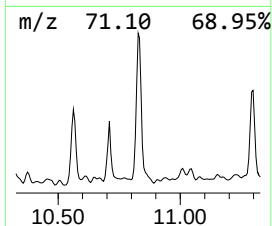
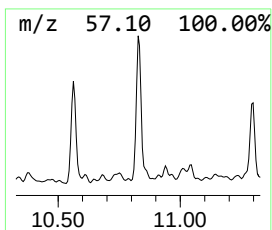
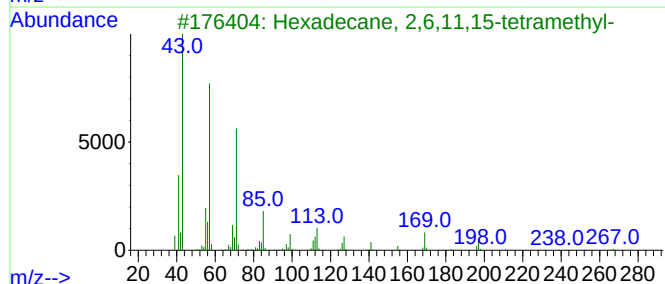
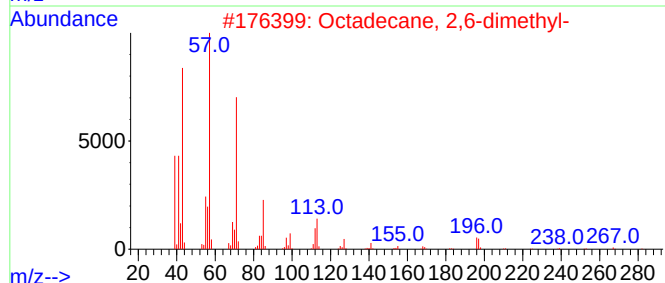
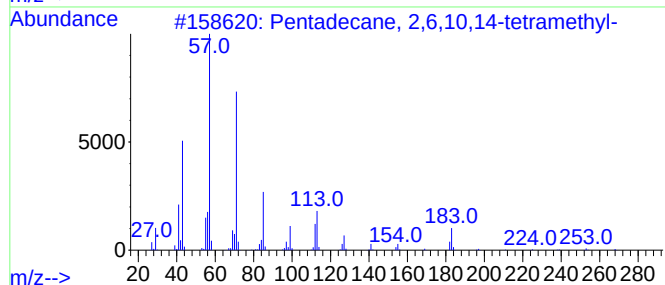
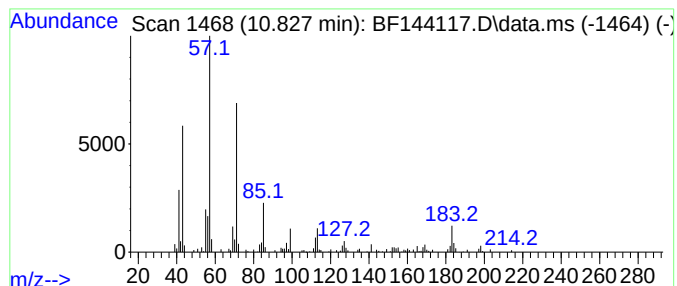
Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Pentadecane, 2,6,10,14-tetr... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD			R.T.
10.827	4.19 ng	117125	Phenanthrene-d10			11.410
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10,14-tetramethyl-		268	C19H40	001921-70-6	86
2	Octadecane, 2,6-dimethyl-		282	C20H42	075163-97-2	72
3	Hexadecane, 2,6,11,15-tetramethyl-		282	C20H42	000504-44-9	72
4	Heptadecane, 2,6-dimethyl-		268	C19H40	054105-67-8	64
5	Tridecane, 4-methyl-		198	C14H30	026730-12-1	59



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Instrument :
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ClientSampleId :
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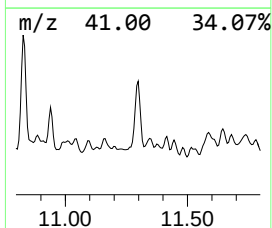
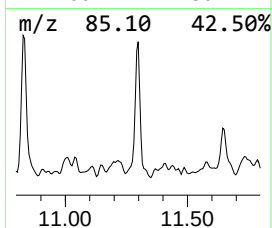
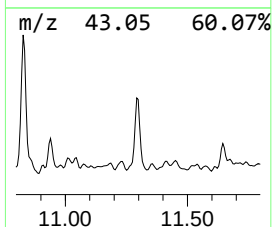
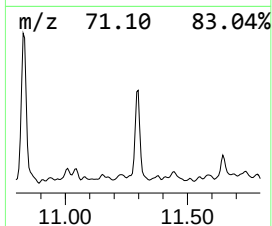
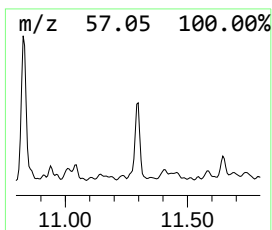
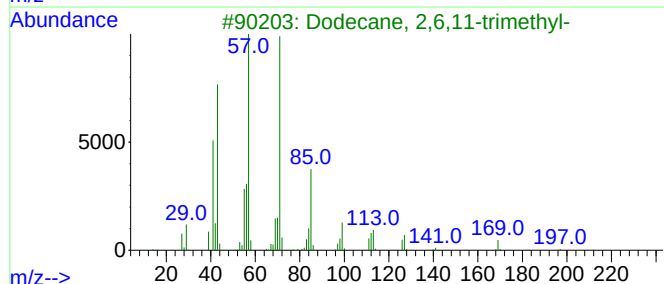
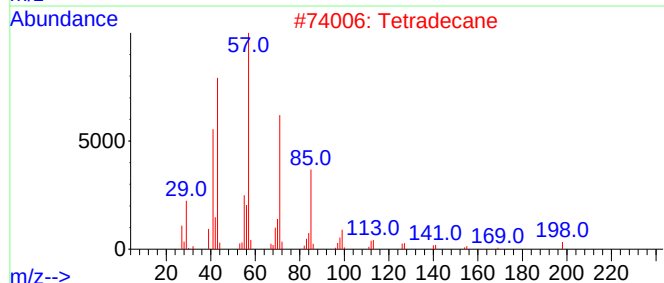
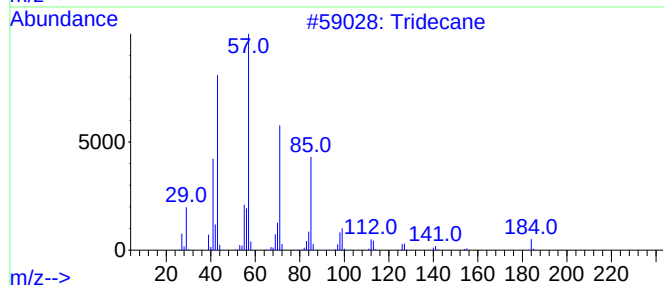
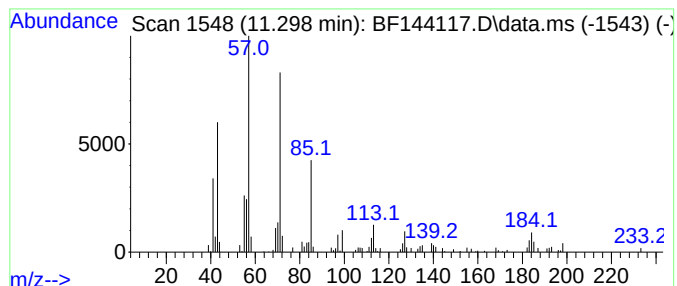
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Tridecane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.298	2.69 ng	75375	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	87
2			Tetradecane	198	C14H30	000629-59-4	87
3			Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	86
4			Tridecane, 3-methyl-	198	C14H30	006418-41-3	80
5			Octane, 2-methyl-	128	C9H20	003221-61-2	74



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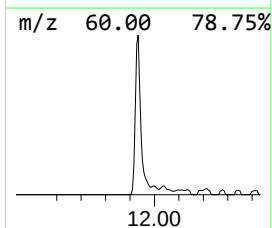
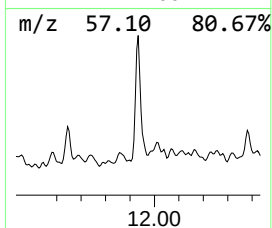
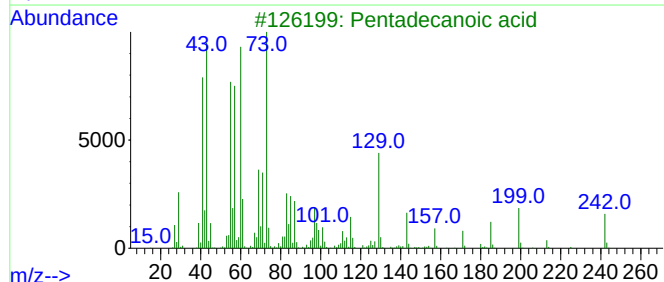
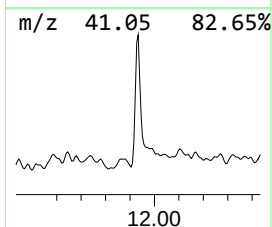
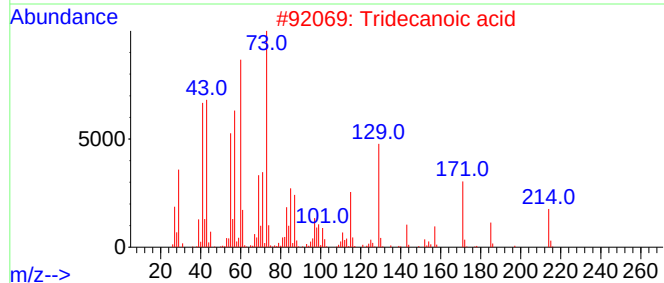
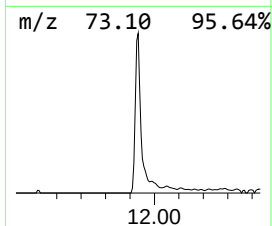
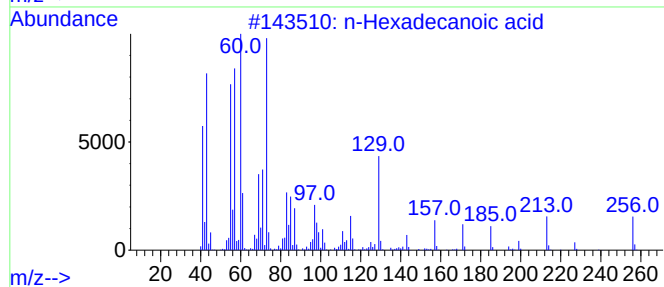
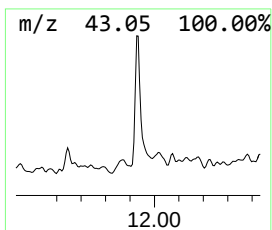
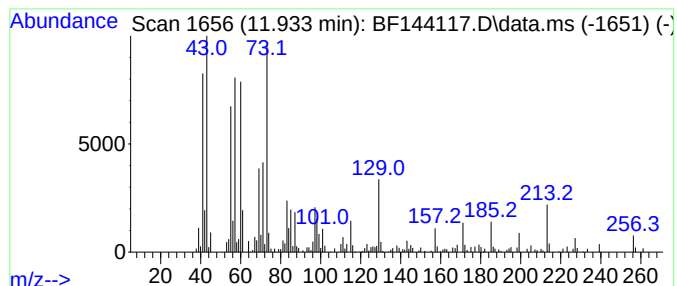
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	6.78 ng	189724	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2			Tridecanoic acid	214	C13H26O2	000638-53-9	92
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4			n-Decanoic acid	172	C10H20O2	000334-48-5	53
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
Acq On : 28 Oct 2025 23:10
Operator : RC/JU
Sample : Q3395-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PR132-S07-102116-20251017

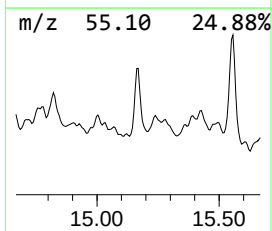
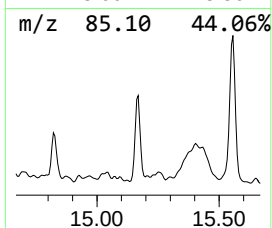
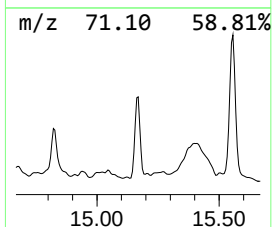
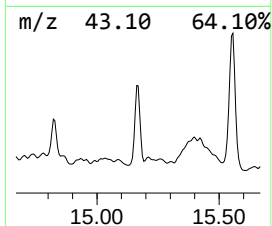
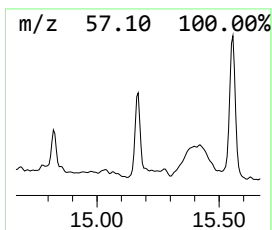
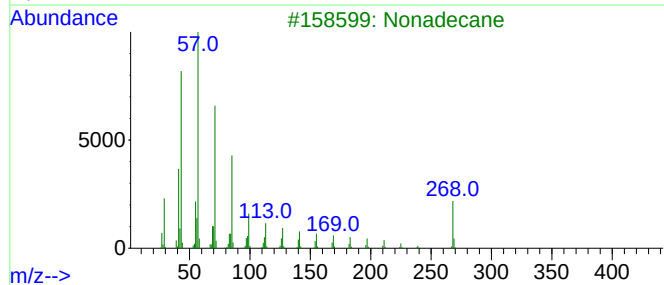
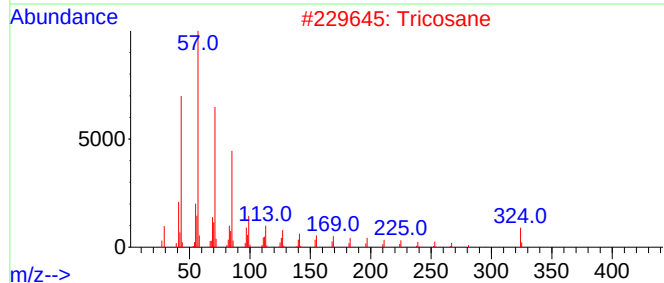
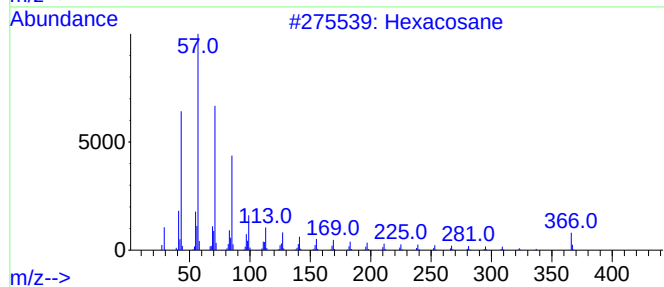
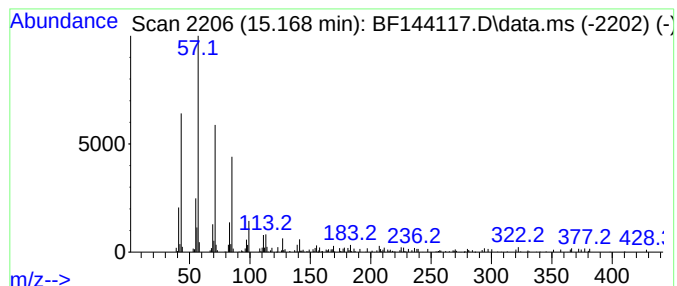
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Hexacosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.168	3.17 ng	103663	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexacosane	366	C26H54	000630-01-3	93
2		Tricosane	324	C23H48	000638-67-5	91
3		Nonadecane	268	C19H40	000629-92-5	87
4		Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	87
5		Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
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Operator : RC/JU
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Instrument :
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ClientSampleId :
PR132-S07-102116-20251017

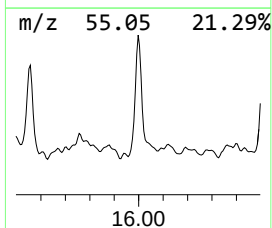
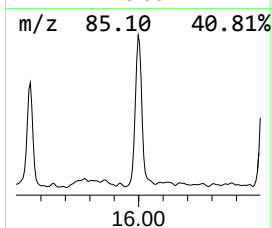
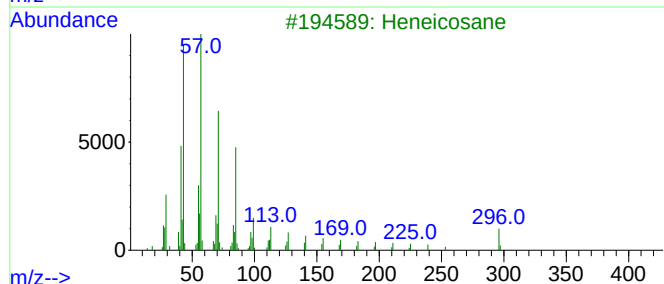
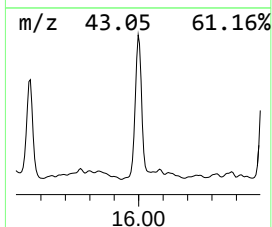
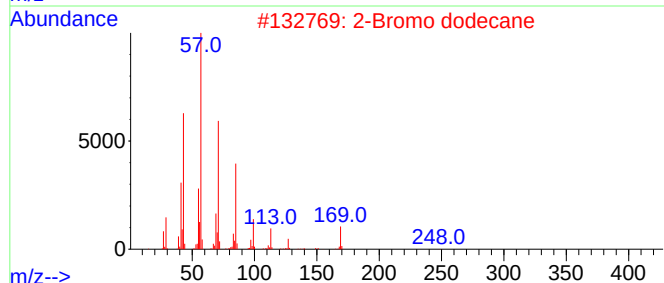
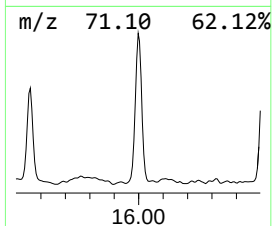
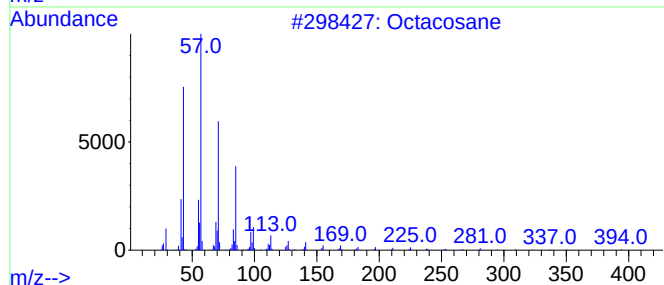
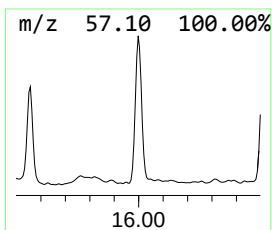
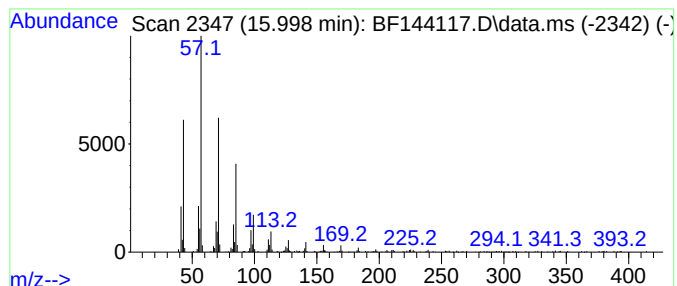
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octacosane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.998	10.10 ng	330190	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	99
2			2-Bromo dodecane	248	C12H25Br	013187-99-0	93
3			Heneicosane	296	C21H44	000629-94-7	91
4			Hexacosane	366	C26H54	000630-01-3	91
5			Hentriacontane	437	C31H64	000630-04-6	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
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Operator : RC/JU
Sample : Q3395-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

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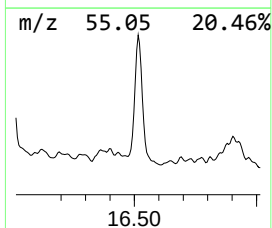
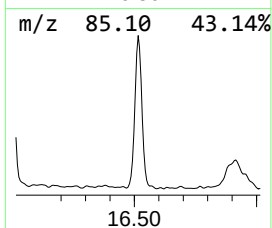
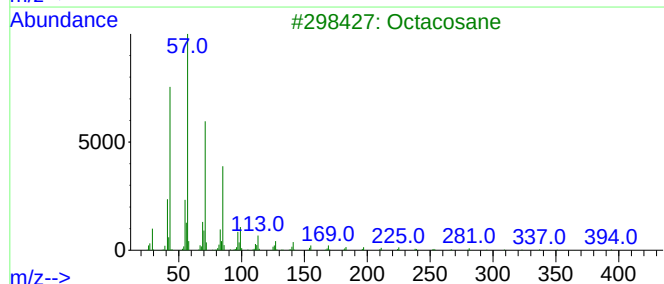
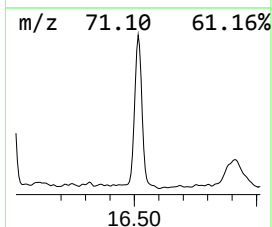
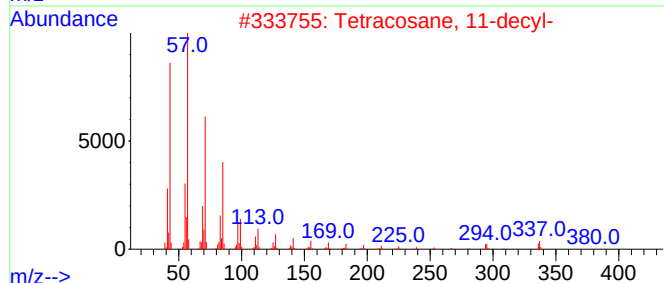
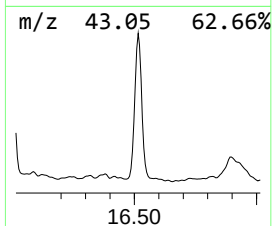
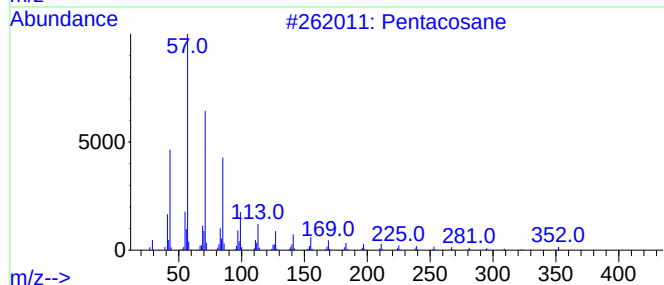
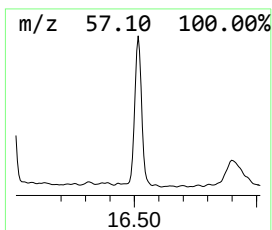
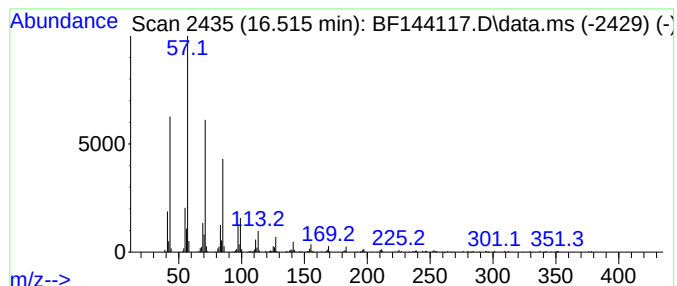
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 Pentacosane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.515	14.15 ng	462449	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	91
2			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Tetratetracontane	619	C44H90	007098-22-8	91
5			Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
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Acq On : 28 Oct 2025 23:10
Operator : RC/JU
Sample : Q3395-05
Misc :
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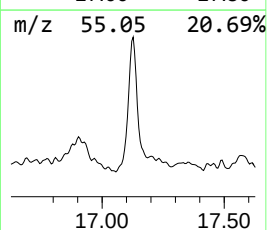
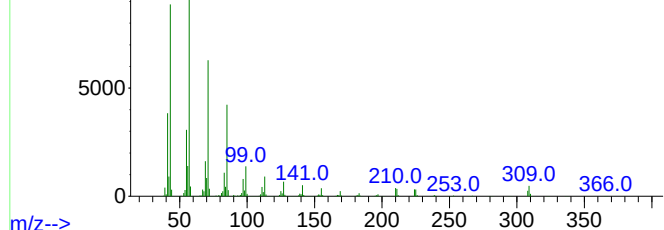
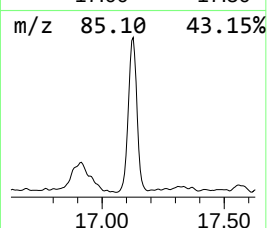
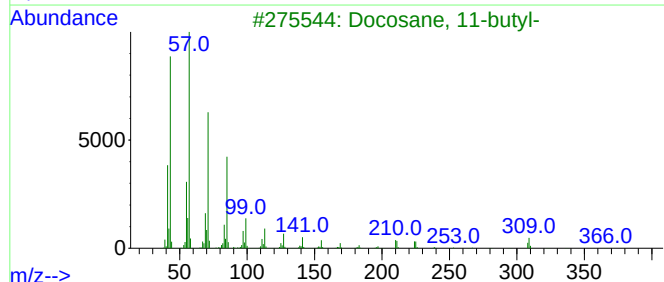
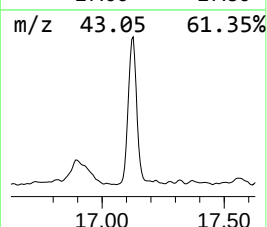
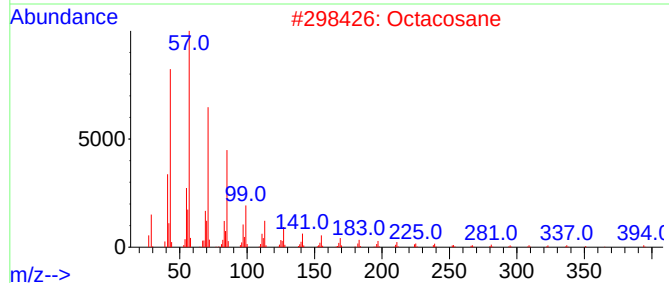
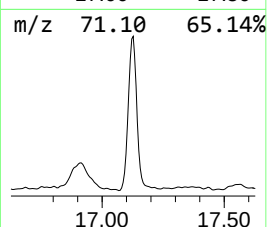
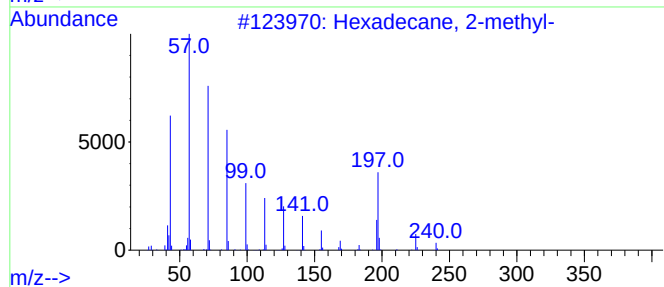
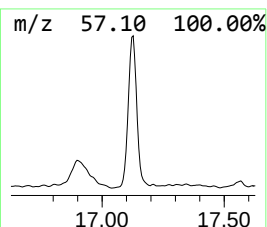
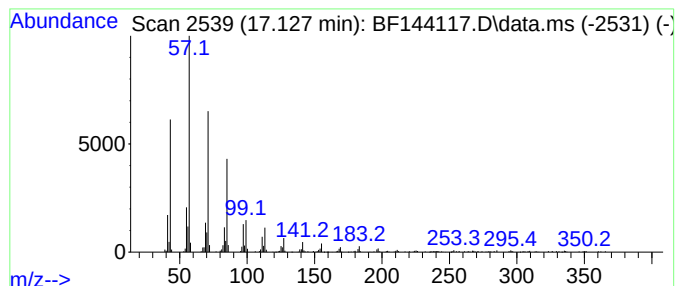
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Hexadecane, 2-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.127	18.36 ng	600053	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	91
2			Octacosane	394	C28H58	000630-02-4	91
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	91
4			Heneicosane, 11-decyl-	437	C31H64	055320-06-4	91
5			4-Methylheneicosane	310	C22H46	025117-29-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
Acq On : 28 Oct 2025 23:10
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PR132-S07-102116-20251017

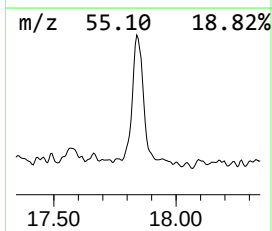
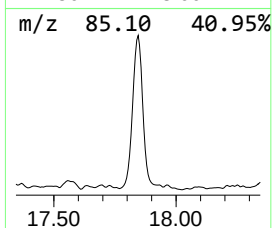
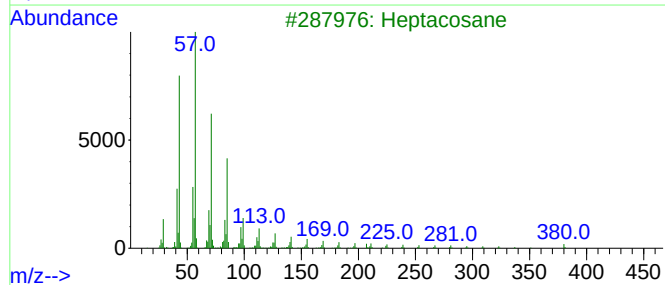
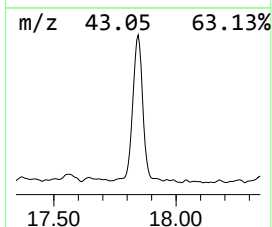
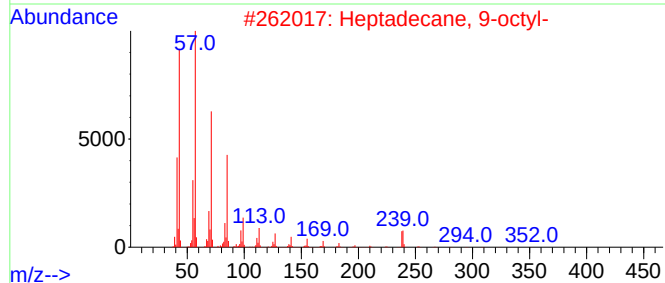
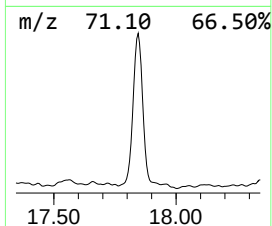
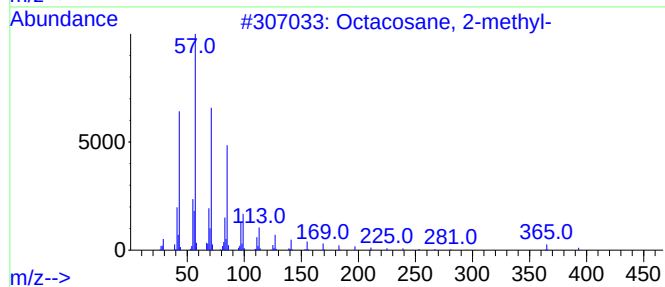
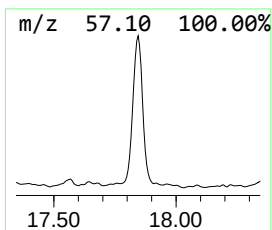
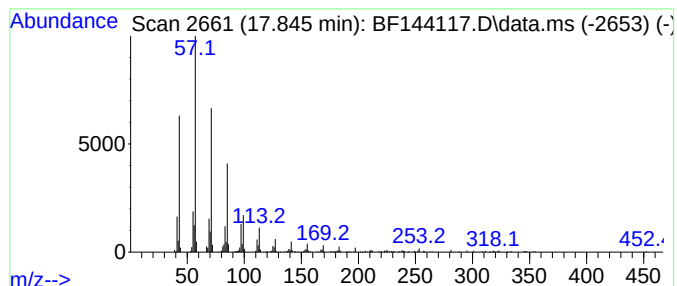
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Octacosane, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.845	17.42 ng	569459	Perylene-d12	15.539

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Hexacosane	366	C26H54	000630-01-3	91
5		Hentriacontane	437	C31H64	000630-04-6	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
Acq On : 28 Oct 2025 23:10
Operator : RC/JU
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ClientSampleId :
PR132-S07-102116-20251017

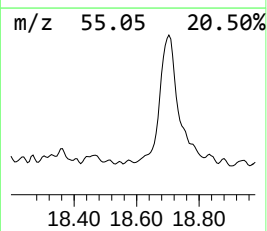
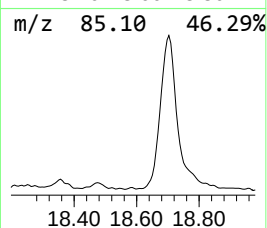
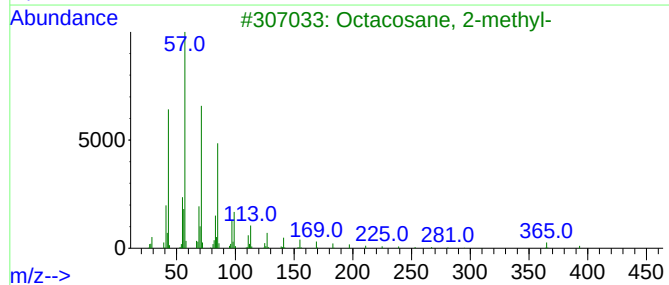
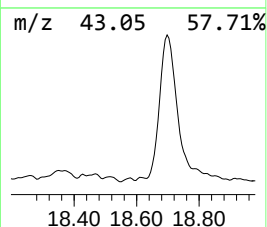
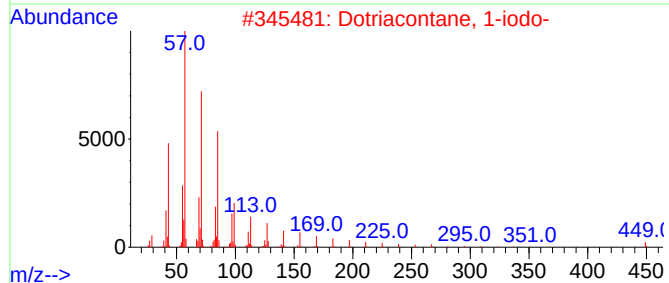
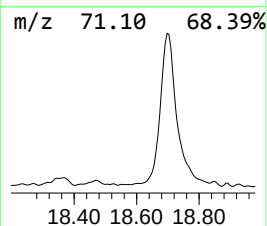
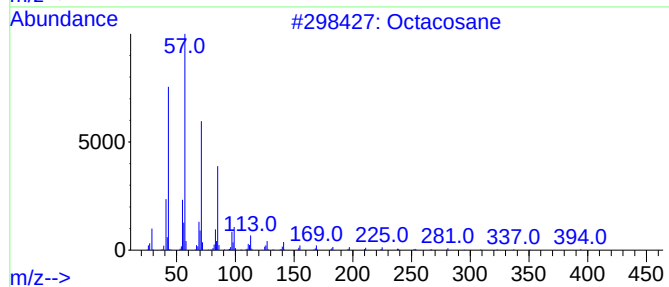
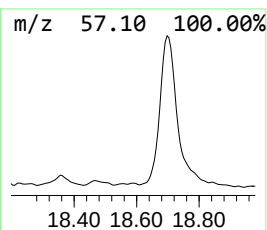
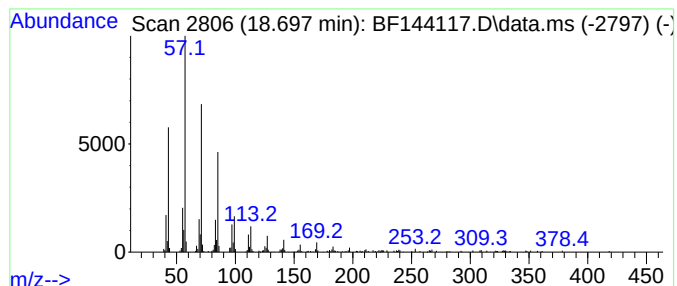
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 15 Dotriacontane, 1-iodo- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.698	19.66 ng	642687	Perylene-d12	15.539

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosane	394	C28H58	000630-02-4	91
2			Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	91
3			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
4			Tetradecane, 2,6,10-trimethyl-	240	C17H36	014905-56-7	90
5			Docosane	310	C22H46	000629-97-0	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102825\
Data File : BF144117.D
Acq On : 28 Oct 2025 23:10
Operator : RC/JU
Sample : Q3395-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PR132-S07-102116-20251017

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF102425.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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					#	RT	Resp	Conc
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Benzophenone	10.633	12.8	ng	370317	3	9.916	578179	20.0
Pentadecane, 2,...	10.827	4.2	ng	117125	4	11.410	559493	20.0
Tridecane	11.298	2.7	ng	75375	4	11.410	559493	20.0
n-Hexadecanoic ...	11.933	6.8	ng	189724	4	11.410	559493	20.0
Hexacosane	15.168	3.2	ng	103663	6	15.539	653749	20.0
Octacosane	15.998	10.1	ng	330190	6	15.539	653749	20.0
Pentacosane	16.515	14.2	ng	462449	6	15.539	653749	20.0
Hexadecane, 2-m...	17.127	18.4	ng	600053	6	15.539	653749	20.0
Octacosane, 2-m...	17.845	17.4	ng	569459	6	15.539	653749	20.0
Dotriacontane, ...	18.698	19.7	ng	642687	6	15.539	653749	20.0