

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
 Data File : BP026335.D
 Acq On : 15 Dec 2025 21:41
 Operator : RC/JU
 Sample : Q3859-01 5X
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HD-02-121225

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-BP111825.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP026335.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.814	314	319	330	rVB	124336	183275	5.20%	0.650%
2	5.278	391	398	408	rBV	470358	735317	20.84%	2.609%
3	6.837	654	663	681	rBV	380968	686773	19.47%	2.437%
4	7.649	794	801	809	rBV	817853	1338402	37.94%	4.749%
5	8.790	979	995	1003	rBV	233910	470631	13.34%	1.670%
6	8.872	1005	1009	1018	rVB	43460	74837	2.12%	0.266%
7	10.413	1261	1271	1283	rBV	1054164	1914696	54.28%	6.794%
8	10.584	1297	1300	1306	rVB	25744	38034	1.08%	0.135%
9	11.125	1380	1392	1395	rBV	21243	67308	1.91%	0.239%
10	11.260	1408	1415	1421	rBV6	20506	37846	1.07%	0.134%
11	11.343	1421	1429	1434	rBV7	22862	49994	1.42%	0.177%
12	11.449	1441	1447	1456	rBV4	41408	98658	2.80%	0.350%
13	11.607	1467	1474	1479	rBV8	20196	39062	1.11%	0.139%
14	11.849	1508	1515	1525	rBV	97688	183116	5.19%	0.650%
15	12.072	1550	1553	1561	rVB5	24030	46248	1.31%	0.164%
16	12.337	1594	1598	1602	rVB5	30999	47795	1.35%	0.170%
17	12.484	1614	1623	1627	rBV6	18383	47308	1.34%	0.168%
18	12.584	1636	1640	1646	rBV4	50809	88296	2.50%	0.313%
19	12.654	1647	1652	1665	rVB2	124393	280046	7.94%	0.994%
20	12.813	1675	1679	1684	rVB3	64560	82175	2.33%	0.292%
21	12.884	1684	1691	1701	rBV	591499	1056575	29.95%	3.749%
22	13.101	1717	1728	1735	rBV3	152245	332377	9.42%	1.179%
23	13.190	1738	1743	1751	rVV5	60380	147326	4.18%	0.523%
24	13.266	1751	1756	1761	rVV	80288	151965	4.31%	0.539%
25	13.331	1762	1767	1777	rVB	60810	156011	4.42%	0.554%
26	13.643	1816	1820	1824	rVB6	24886	40521	1.15%	0.144%
27	13.695	1825	1829	1833	rVB6	25448	38665	1.10%	0.137%
28	13.743	1833	1837	1838	rBV3	32654	41770	1.18%	0.148%
29	13.772	1839	1842	1845	rVV2	56931	68167	1.93%	0.242%
30	13.995	1875	1880	1884	rBV5	18898	41190	1.17%	0.146%
31	14.119	1898	1901	1907	rVB5	25453	50507	1.43%	0.179%
32	14.184	1908	1912	1917	rBV	201227	267142	7.57%	0.948%
33	14.272	1920	1927	1935	rBV	1607368	2408415	68.27%	8.546%
34	14.666	1988	1994	1997	rBV7	22918	46712	1.32%	0.166%
35	14.825	2017	2021	2027	rVB2	71566	112116	3.18%	0.398%
36	14.895	2029	2033	2036	rBV3	32845	44948	1.27%	0.159%

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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_P\Methods\8270E-BP111825.M
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37	15.166	2074	2079	2083	rBV	232611	289552	8.21%	1.027%
38	15.342	2106	2109	2114	rBV5	21333	35336	1.00%	0.125%
39	15.578	2142	2149	2154	rBV	100145	189078	5.36%	0.671%
40	15.660	2159	2163	2170	rVB2	82212	156799	4.44%	0.556%
41	15.737	2172	2176	2180	rBV4	33343	57398	1.63%	0.204%
42	15.795	2181	2186	2195	rBV	556827	891426	25.27%	3.163%
43	16.042	2223	2228	2231	rBV	241210	360857	10.23%	1.280%
44	16.866	2364	2368	2372	rBV	229955	289470	8.21%	1.027%
45	16.919	2373	2377	2382	rVB	125681	182836	5.18%	0.649%
46	17.084	2399	2405	2410	rBV2	1609392	2617600	74.20%	9.288%
47	17.125	2410	2412	2418	rVB	133431	179607	5.09%	0.637%
48	17.554	2481	2485	2489	rBV4	35886	52089	1.48%	0.185%
49	17.625	2493	2497	2506	rVB	225776	281766	7.99%	1.000%
50	17.813	2525	2529	2532	rBV	62487	80192	2.27%	0.285%
51	18.036	2563	2567	2575	rVB2	204285	303902	8.61%	1.078%
52	18.195	2591	2594	2599	rVB4	58461	83968	2.38%	0.298%
53	18.348	2616	2620	2627	rBV	185564	257709	7.31%	0.914%
54	19.019	2727	2734	2740	rBV5	185315	470641	13.34%	1.670%
55	19.225	2765	2769	2781	rBV	360600	787324	22.32%	2.794%
56	19.378	2791	2795	2804	rBV4	187873	316698	8.98%	1.124%
57	19.607	2830	2834	2843	rVB2	318897	602245	17.07%	2.137%
58	19.830	2867	2872	2878	rBV	944189	1448131	41.05%	5.139%
59	20.183	2929	2932	2935	rBV	136289	167096	4.74%	0.593%
60	21.530	3155	3161	3166	rBV2	1762883	3037439	86.10%	10.778%
61	24.836	3715	3723	3733	rVB	1261771	3527714	100.00%	12.518%

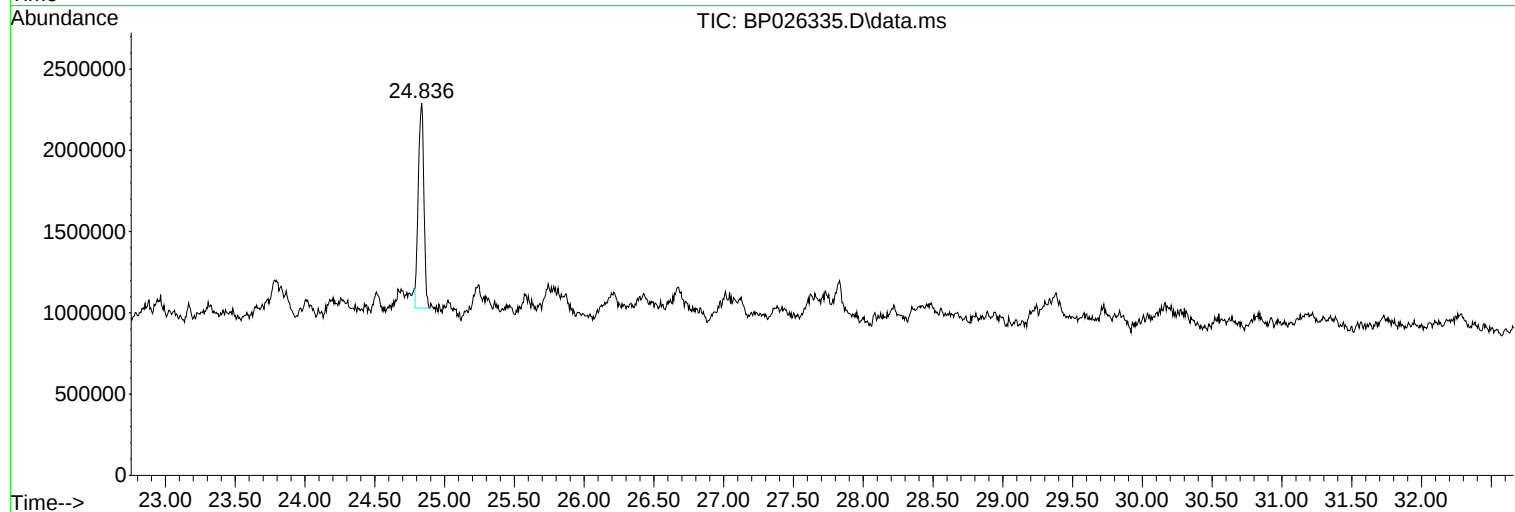
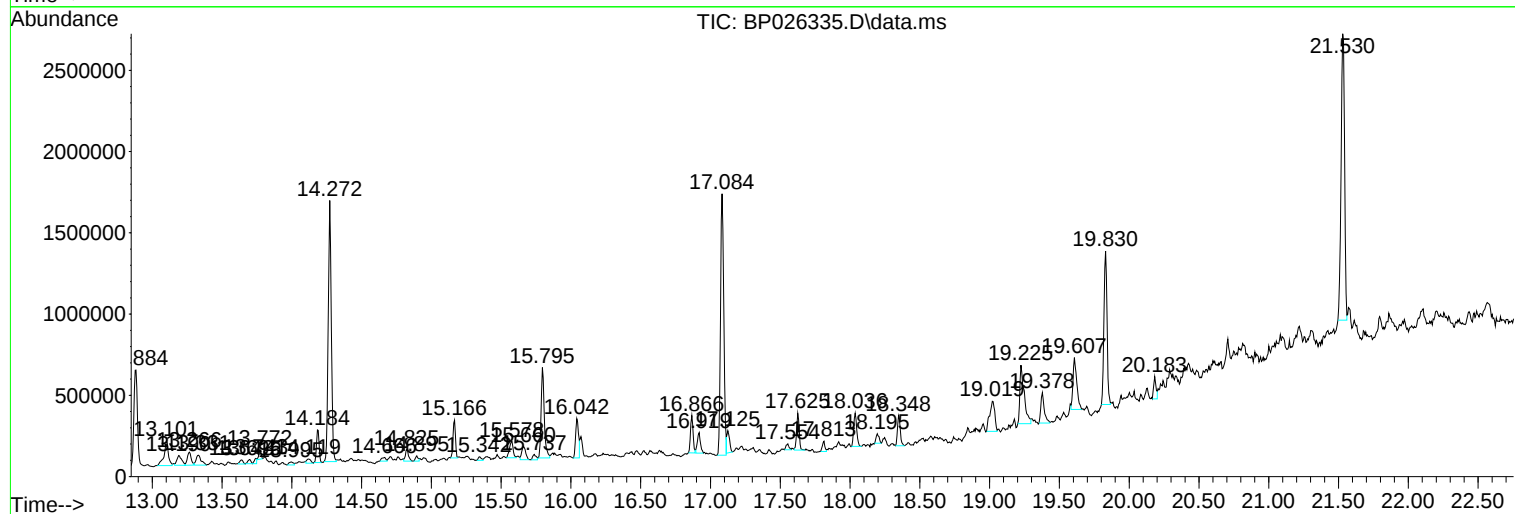
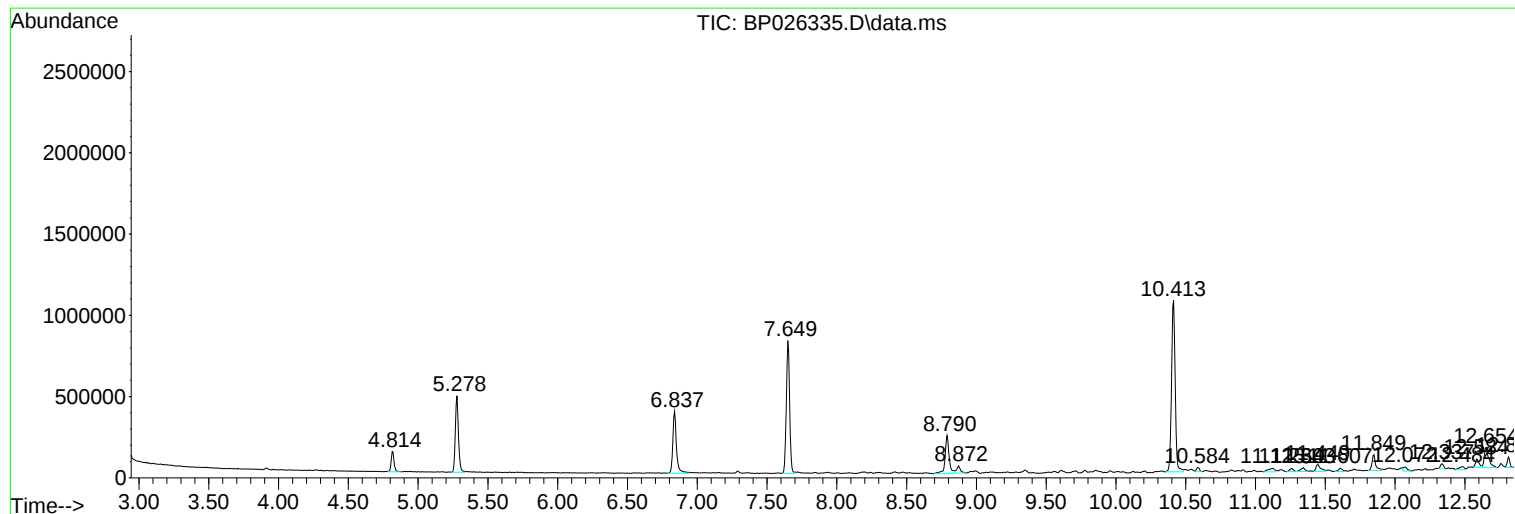
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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-02-121225

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.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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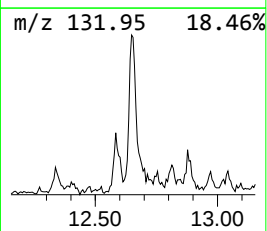
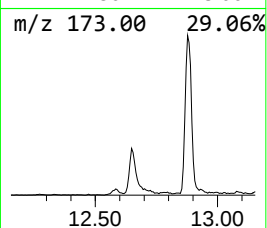
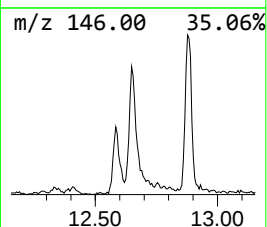
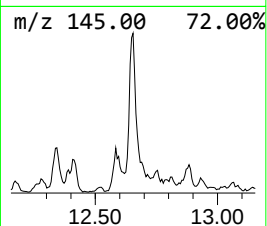
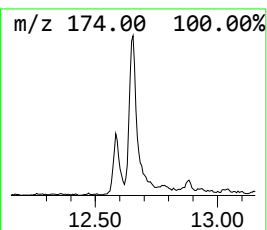
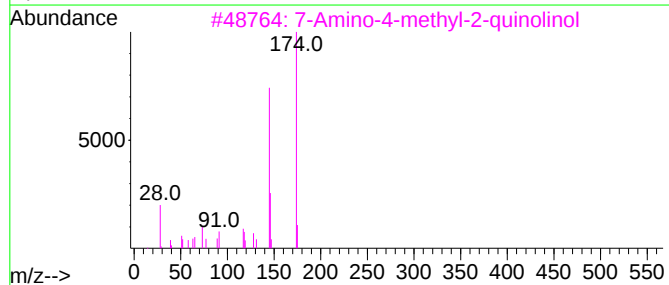
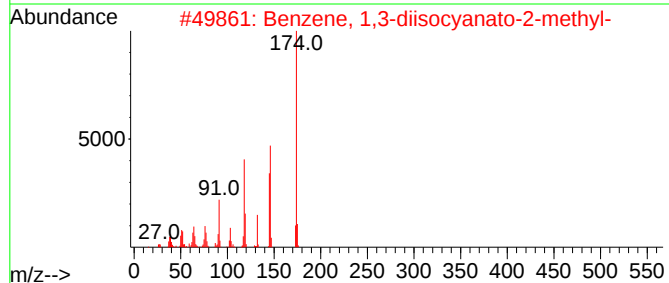
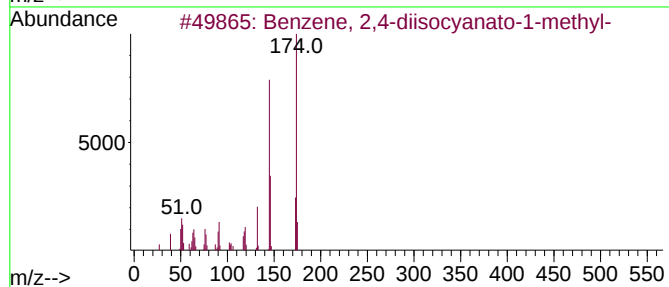
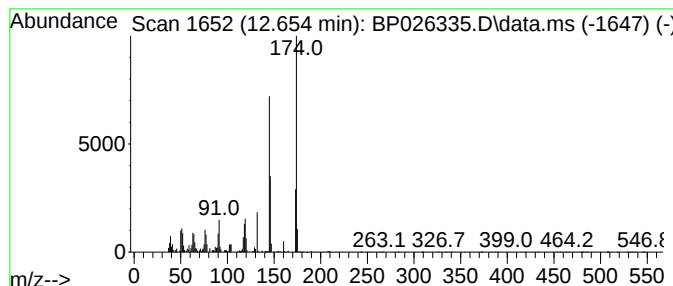
Instrument :
BNA_P
ClientSampleId :
HD-02-121225

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 Benzene, 2,4-diisocyanato-1... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD		R.T.
12.654	2.33 ng	280046	Acenaphthene-d10		14.272
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2,4-diisocyanato-1-methyl-	174	C9H6N2O2	000584-84-9	97
2	Benzene, 1,3-diisocyanato-2-methyl-	174	C9H6N2O2	000091-08-7	64
3	7-Amino-4-methyl-2-quinolinol	174	C10H10N2O	019840-99-4	64
4	Benzo [f]-1,5-diazabicyclo[3.2.2...]	174	C11H14N2	007092-76-4	64
5	3-Ethyl-1,2-dihydro-2-oxoquinox...	174	C10H10N2O	013297-35-3	53



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BNA_P
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HD-02-121225

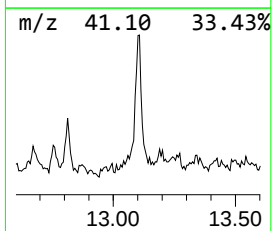
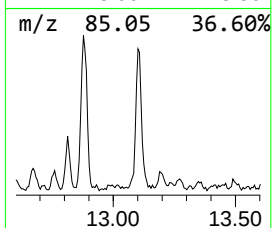
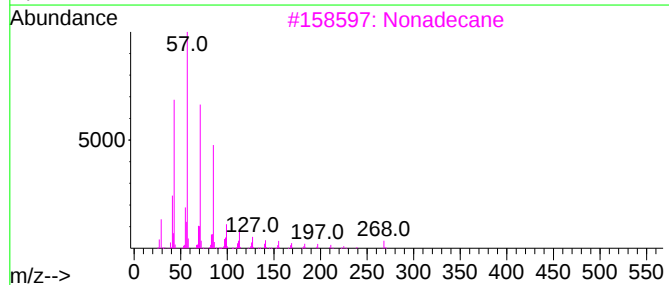
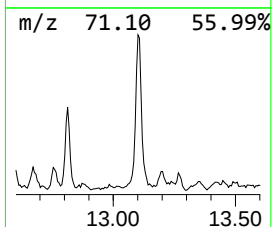
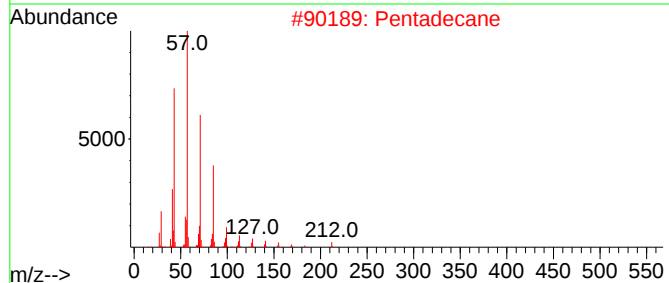
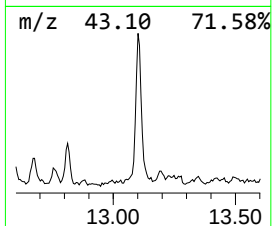
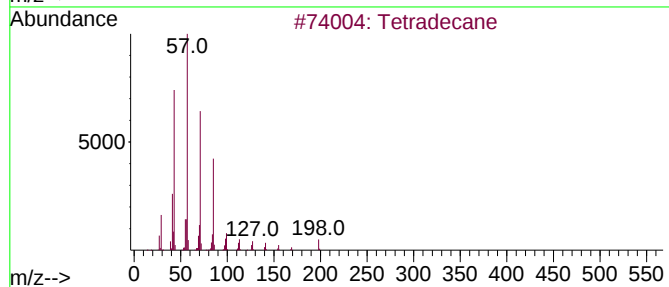
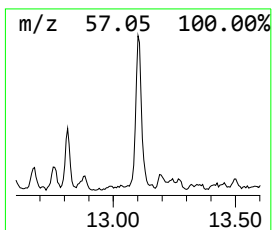
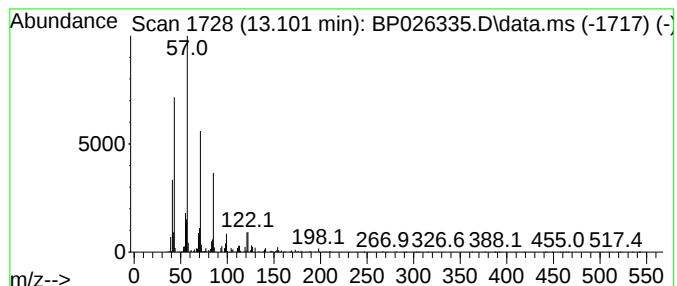
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 3 Tetradecane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.101	2.76 ng	332377	Acenaphthene-d10	14.272

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetradecane	198	C14H30	000629-59-4	95
2			Pentadecane	212	C15H32	000629-62-9	81
3			Nonadecane	268	C19H40	000629-92-5	81
4			Nonadecane, 9-methyl-	282	C20H42	013287-24-6	81
5			Eicosane	282	C20H42	000112-95-8	76



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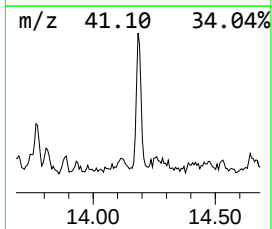
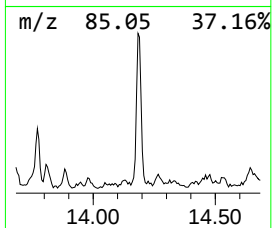
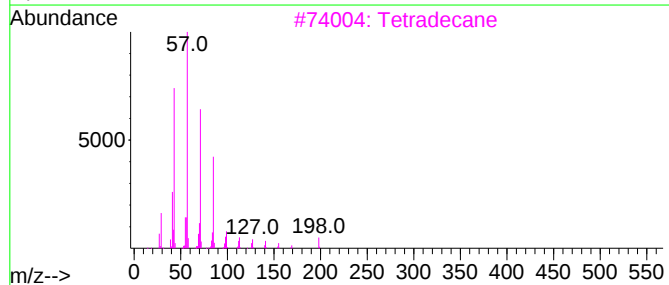
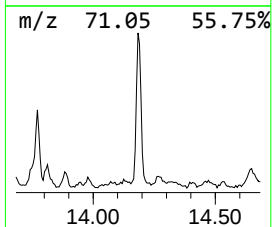
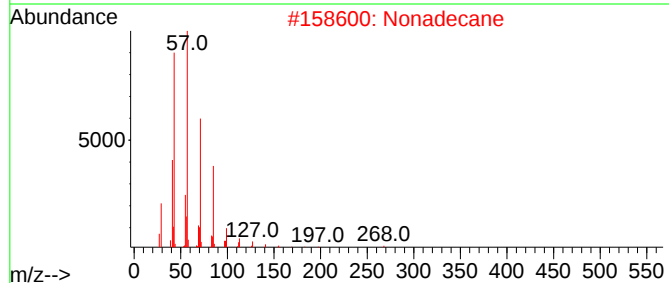
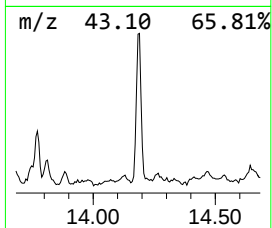
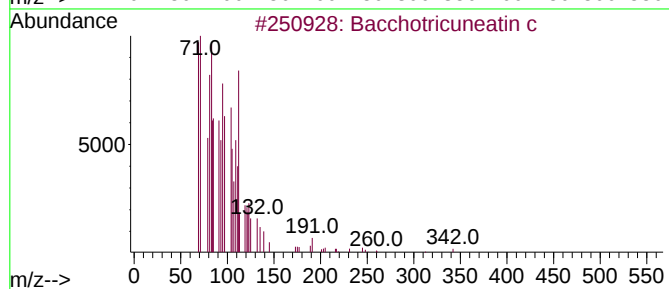
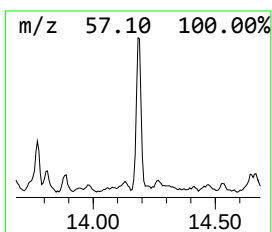
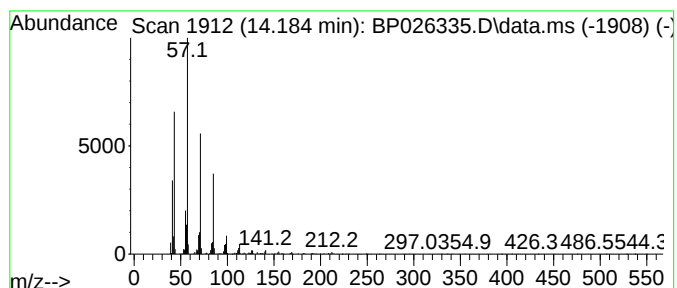
Instrument :
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Bacchotricuneatin c Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.184	2.22 ng	267142	Acenaphthene-d10		14.272	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Bacchotricuneatin c		342	C20H22O5	066563-30-2	95
2	Nonadecane		268	C19H40	000629-92-5	91
3	Tetradecane		198	C14H30	000629-59-4	90
4	Tridecane, 6-methyl-		198	C14H30	013287-21-3	90
5	Hexadecane		226	C16H34	000544-76-3	90



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Instrument :
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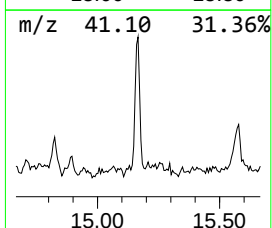
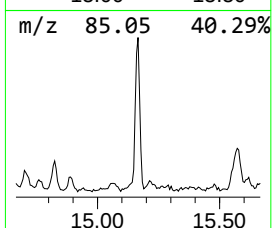
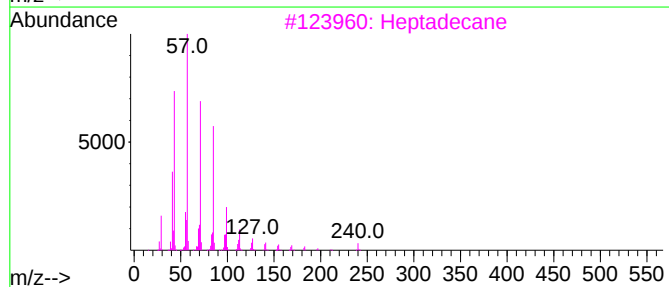
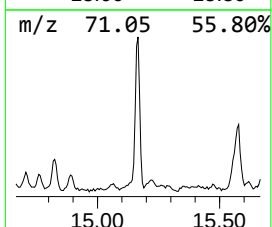
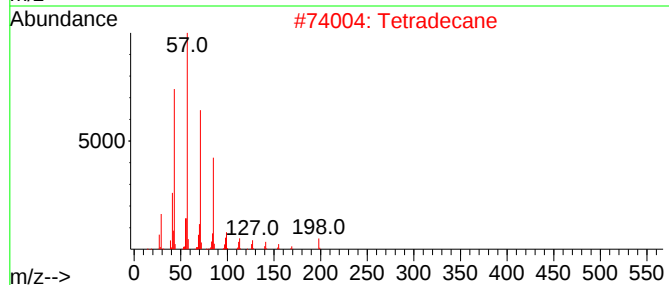
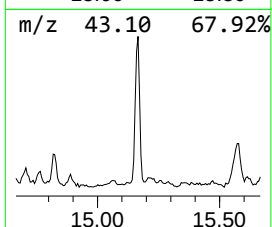
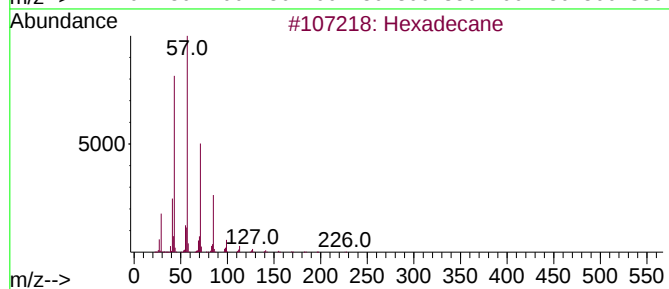
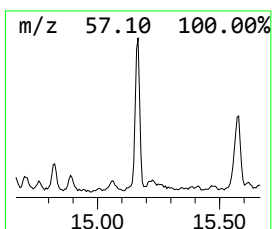
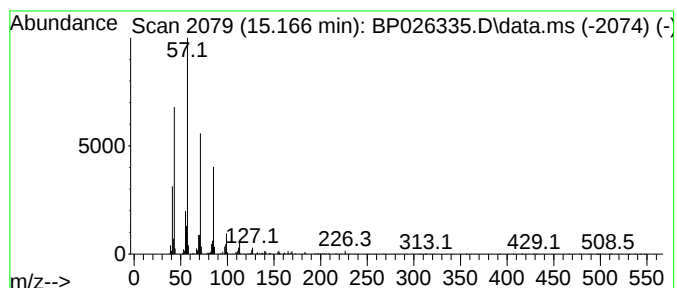
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP111825.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Hexadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.166	2.40 ng	289552	Acenaphthene-d10	14.272

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	97
2			Tetradecane	198	C14H30	000629-59-4	93
3			Heptadecane	240	C17H36	000629-78-7	90
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	87
5			Nonadecane	268	C19H40	000629-92-5	83



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ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-02-121225

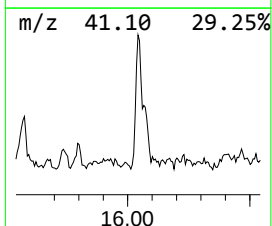
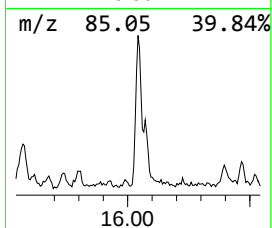
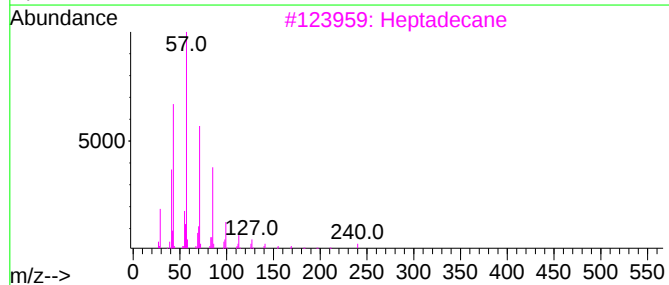
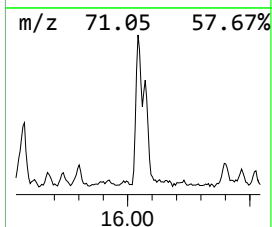
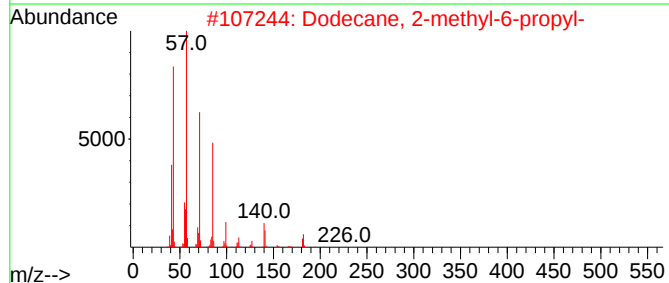
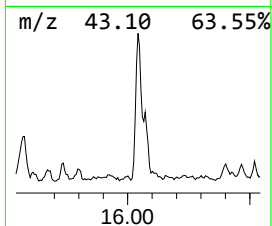
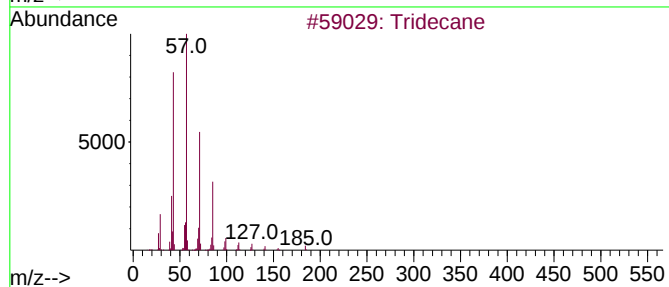
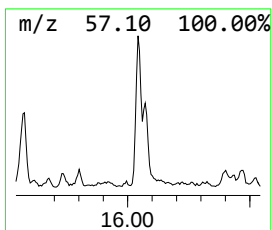
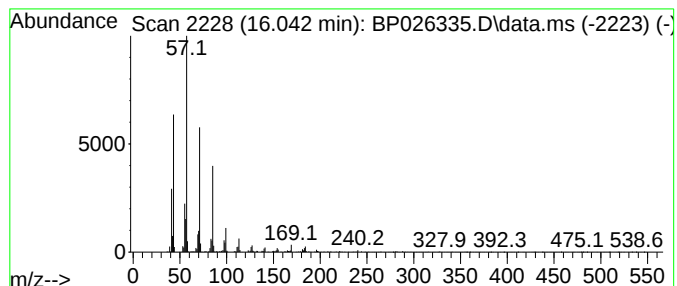
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Tridecane Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.042	2.76 ng	360857	Phenanthrene-d10	17.084

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	93
2			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
3			Heptadecane	240	C17H36	000629-78-7	91
4			Nonadecane	268	C19H40	000629-92-5	91
5			Heptacosane	380	C27H56	000593-49-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
Data File : BP026335.D
Acq On : 15 Dec 2025 21:41
Operator : RC/JU
Sample : Q3859-01 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-02-121225

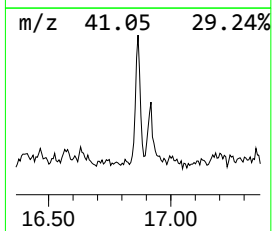
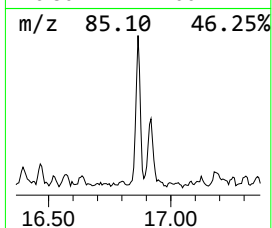
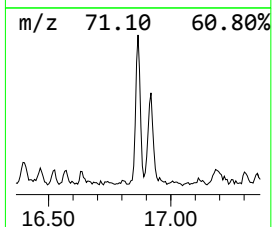
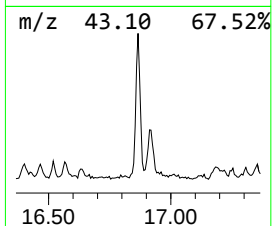
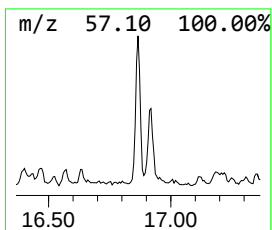
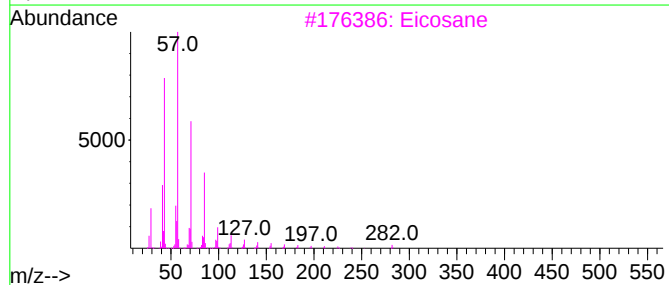
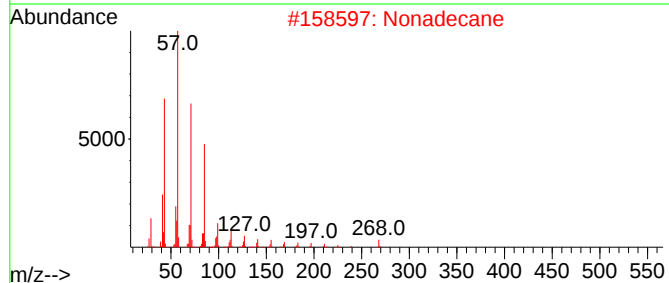
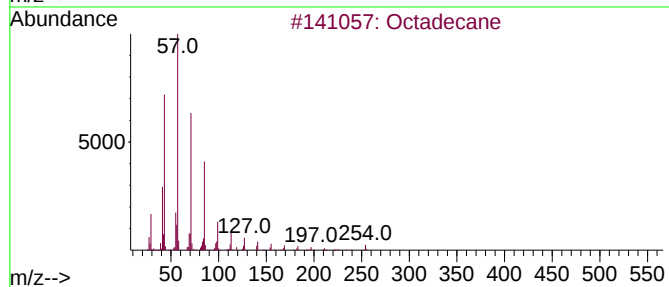
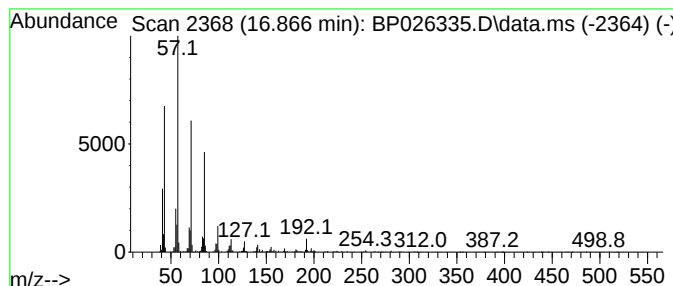
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Octadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.866	2.21 ng	289470	Phenanthrene-d10	17.084

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	93
2			Nonadecane	268	C19H40	000629-92-5	91
3			Eicosane	282	C20H42	000112-95-8	87
4			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
5			Tetradecane	198	C14H30	000629-59-4	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
Data File : BP026335.D
Acq On : 15 Dec 2025 21:41
Operator : RC/JU
Sample : Q3859-01 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HD-02-121225

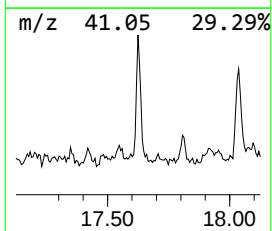
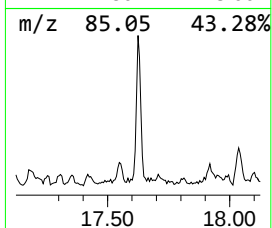
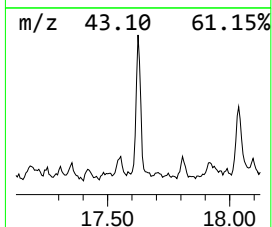
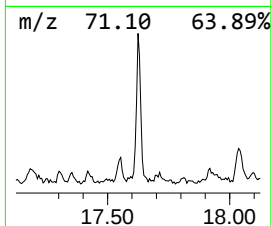
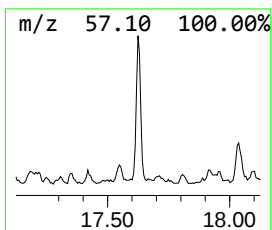
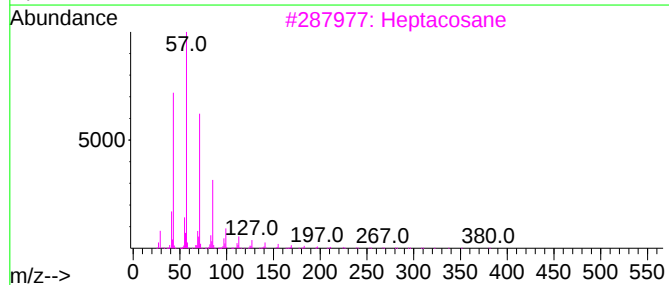
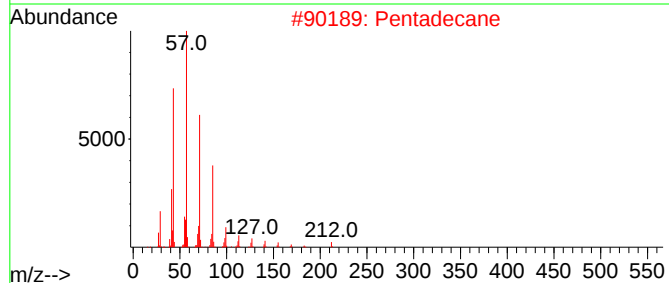
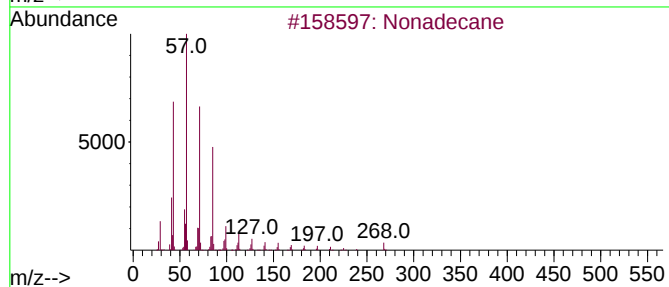
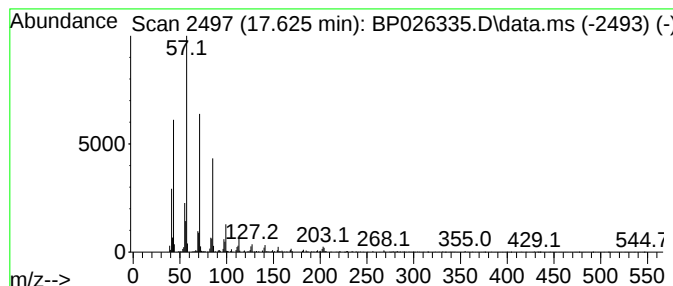
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Nonadecane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.625	2.15 ng	281766	Phenanthrene-d10	17.084

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	97
2			Pentadecane	212	C15H32	000629-62-9	93
3			Heptacosane	380	C27H56	000593-49-7	91
4			Tetracosane	338	C24H50	000646-31-1	90
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
Data File : BP026335.D
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Operator : RC/JU
Sample : Q3859-01 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

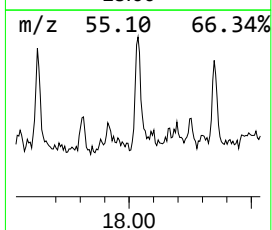
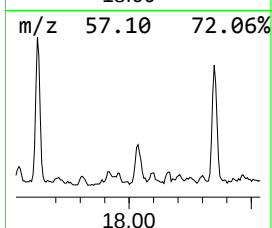
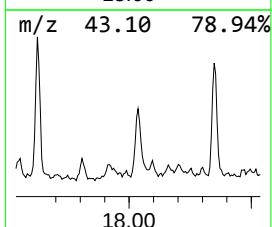
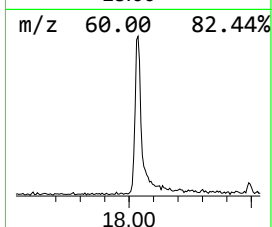
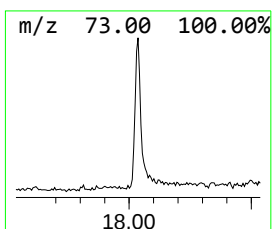
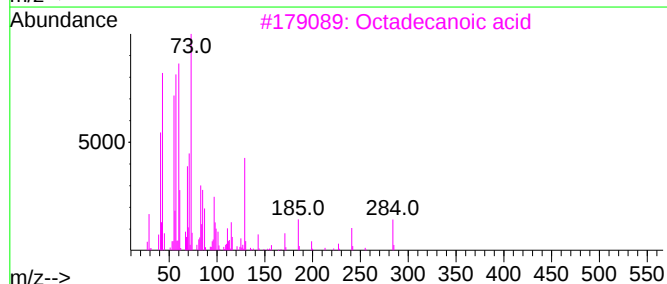
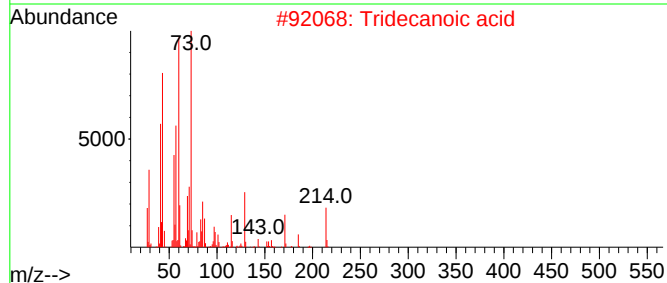
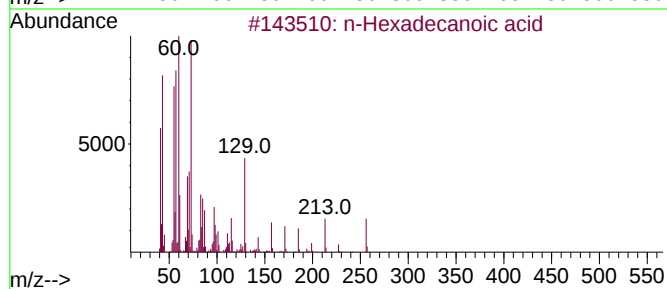
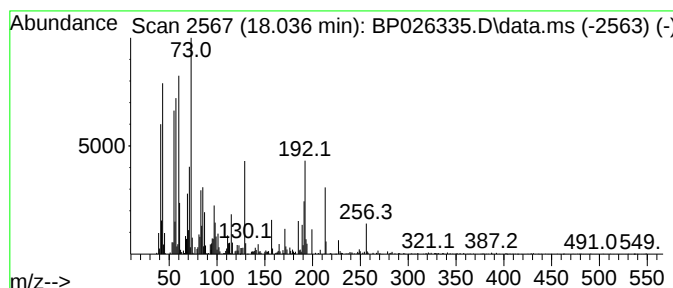
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.036	2.32 ng	303902	Phenanthrene-d10	17.084	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2	Tridecanoic acid	214	C13H26O2	000638-53-9	83
3	Octadecanoic acid	284	C18H36O2	000057-11-4	76
4	Pentadecanoic acid	242	C15H30O2	001002-84-2	76
5	n-Decanoic acid	172	C10H20O2	000334-48-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
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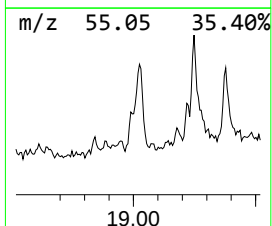
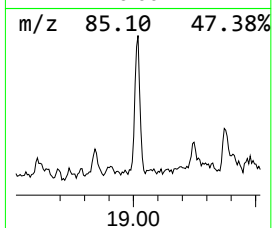
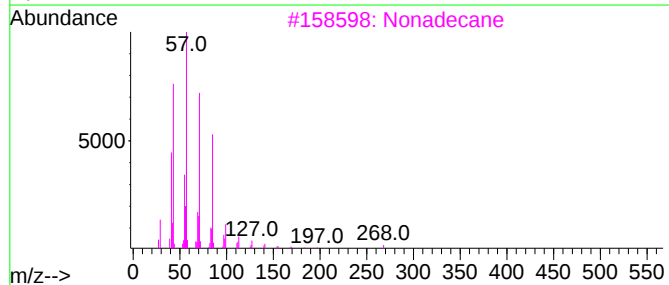
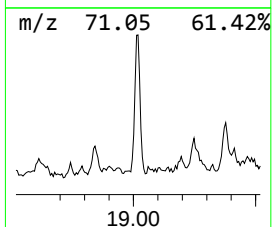
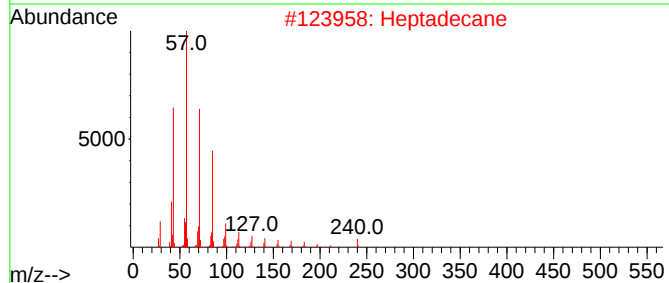
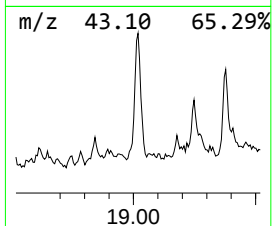
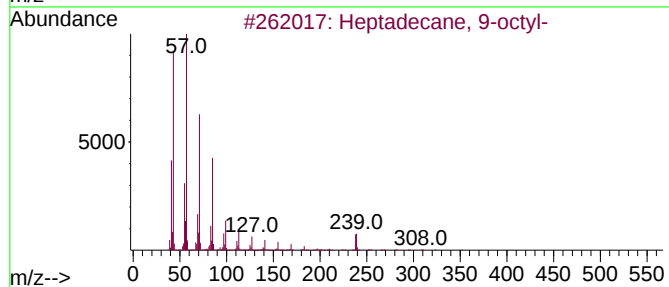
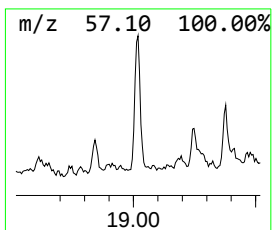
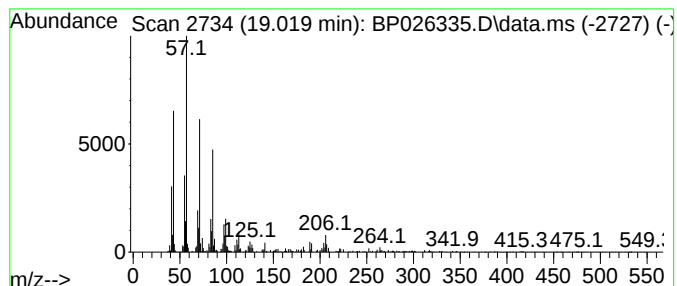
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Heptadecane, 9-octyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.019	3.60 ng	470641	Phenanthrene-d10		17.084	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
2	Heptadecane		240	C17H36	000629-78-7	90
3	Nonadecane		268	C19H40	000629-92-5	90
4	10-Methylnonadecane		282	C20H42	056862-62-5	90
5	Tetratetracontane		619	C44H90	007098-22-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
Data File : BP026335.D
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Sample : Q3859-01 5X
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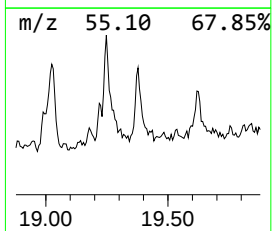
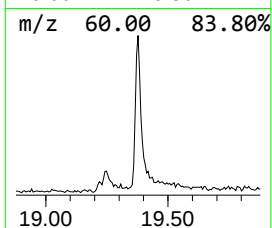
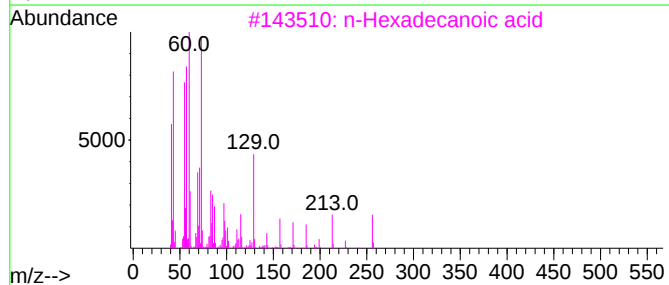
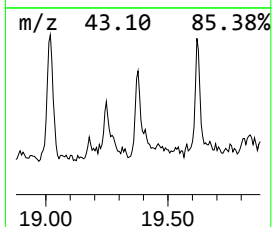
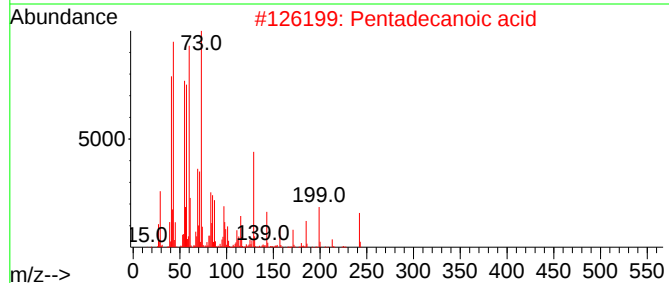
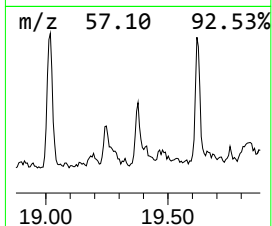
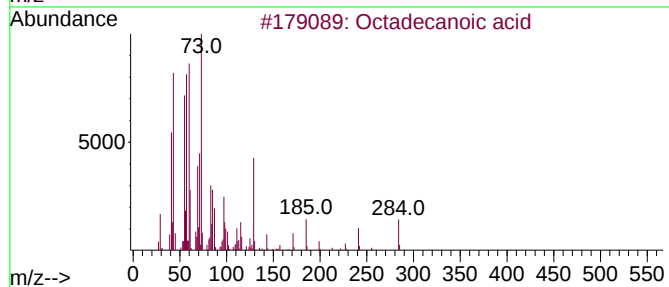
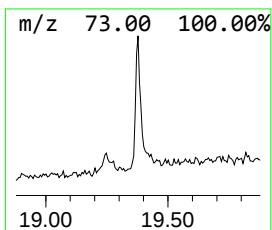
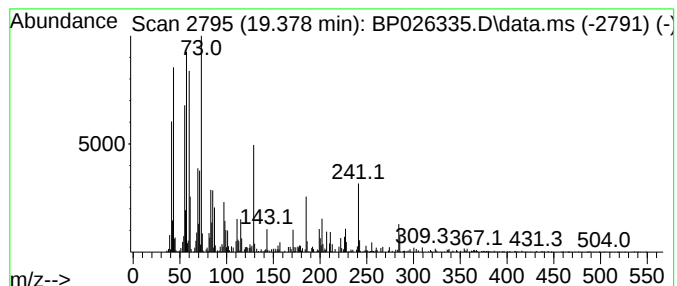
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Octadecanoic acid Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.378	2.09 ng	316698	Chrysene-d12	21.530	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Pentadecanoic acid	242	C15H30O2	001002-84-2	94
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	83
5	Undecanoic acid	186	C11H22O2	000112-37-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121525\
Data File : BP026335.D
Acq On : 15 Dec 2025 21:41
Operator : RC/JU
Sample : Q3859-01 5X
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
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Tetradecane	13.101	2.8	ng	332377	3	14.272	2408420	20.0
Bacchotricuneat...	14.184	2.2	ng	267142	3	14.272	2408420	20.0
Hexadecane	15.166	2.4	ng	289552	3	14.272	2408420	20.0
Tridecane	16.042	2.8	ng	360857	4	17.084	2617600	20.0
Octadecane	16.866	2.2	ng	289470	4	17.084	2617600	20.0
Nonadecane	17.625	2.1	ng	281766	4	17.084	2617600	20.0
n-Hexadecanoic ...	18.036	2.3	ng	303902	4	17.084	2617600	20.0
Heptadecane, 9-...	19.019	3.6	ng	470641	4	17.084	2617600	20.0
Octadecanoic acid	19.378	2.1	ng	316698	5	21.530	3037440	20.0