

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
 Data File : BF138558.D
 Acq On : 12 Jul 2024 21:09
 Operator : CG\JU
 Sample : P3187-09
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-3

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-BF070924.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138558.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.140	3	9	14	rBV	2058495	2557377	100.00%	11.264%
2	5.081	505	509	518	rVB	181448	209664	8.20%	0.923%
3	5.493	575	579	590	rBV	2872329	2467945	96.50%	10.870%
4	6.504	746	751	754	rBV	2584160	2401723	93.91%	10.578%
5	6.869	809	813	816	rBV	779833	698299	27.31%	3.076%
6	7.357	890	896	902	rBV2	24201	49957	1.95%	0.220%
7	7.428	904	908	911	rBV	1817585	1475732	57.70%	6.500%
8	8.145	1027	1030	1034	rBV	966185	779585	30.48%	3.434%
9	9.222	1209	1213	1216	rBV	2853225	2193661	85.78%	9.662%
10	9.763	1302	1305	1308	rBV	98716	81904	3.20%	0.361%
11	9.898	1325	1328	1332	rBV	885295	727173	28.43%	3.203%
12	9.992	1339	1344	1349	rBV	60338	80599	3.15%	0.355%
13	10.686	1458	1462	1466	rBV	1365735	1201493	46.98%	5.292%
14	10.816	1479	1484	1489	rBV3	65797	79047	3.09%	0.348%
15	10.869	1490	1493	1495	rBV	32138	31832	1.24%	0.140%
16	10.898	1495	1498	1502	rVB3	29612	43314	1.69%	0.191%
17	11.045	1520	1523	1527	rVB3	44543	42864	1.68%	0.189%
18	11.081	1527	1529	1534	rBV	36846	40990	1.60%	0.181%
19	11.192	1545	1548	1551	rVB	36379	31111	1.22%	0.137%
20	11.333	1568	1572	1575	rBV	109105	100528	3.93%	0.443%
21	11.386	1577	1581	1583	rVV	925259	759631	29.70%	3.346%
22	11.410	1583	1585	1588	rVV	212593	161143	6.30%	0.710%
23	11.463	1591	1594	1596	rBV	126439	115463	4.51%	0.509%
24	11.616	1615	1620	1625	rBV	119566	124213	4.86%	0.547%
25	11.886	1661	1666	1667	rBV3	40850	59156	2.31%	0.261%
26	11.910	1667	1670	1674	rVV2	442451	401776	15.71%	1.770%
27	11.998	1682	1685	1692	rVB2	82873	117439	4.59%	0.517%
28	12.204	1718	1720	1724	rVB	101041	86694	3.39%	0.382%
29	12.445	1757	1761	1762	rBV3	27945	37051	1.45%	0.163%
30	12.469	1762	1765	1768	rVV4	32386	50062	1.96%	0.220%
31	12.522	1771	1774	1777	rVV	75219	70705	2.76%	0.311%
32	12.598	1784	1787	1790	rBV	700139	546744	21.38%	2.408%
33	12.692	1800	1803	1806	rBV	153171	158772	6.21%	0.699%
34	12.827	1820	1826	1829	rBV	706135	661969	25.88%	2.916%
35	12.969	1846	1850	1853	rBV	2341912	1915283	74.89%	8.436%
36	13.045	1860	1863	1867	rBV2	48260	62060	2.43%	0.273%

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Instrument :
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ClientSampleId :
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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-BF070924.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.663	1965	1968	1972	rBV2	88505	130316	5.10%	0.574%
38	13.869	2000	2003	2007	rBV	185545	203640	7.96%	0.897%
39	14.022	2025	2029	2032	rBV2	524111	610491	23.87%	2.689%
40	15.063	2202	2206	2209	rBV	374469	553858	21.66%	2.439%
41	15.363	2255	2257	2263	rVB	182399	209329	8.19%	0.922%
42	15.492	2275	2279	2282	rBV2	289754	373742	14.61%	1.646%

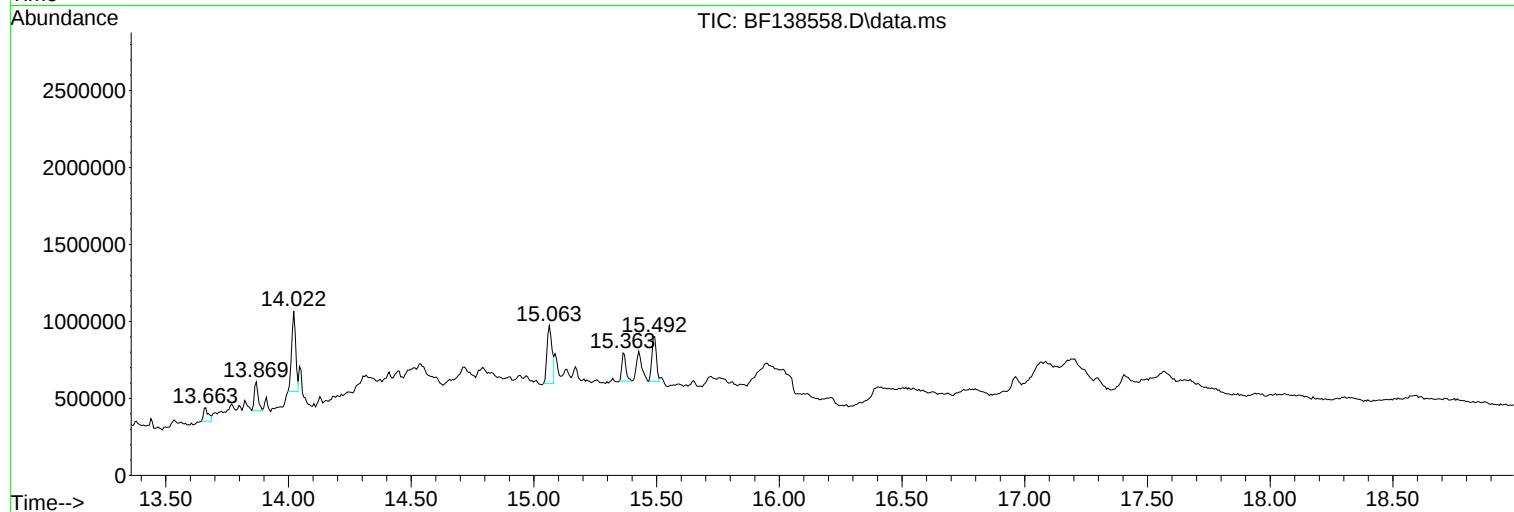
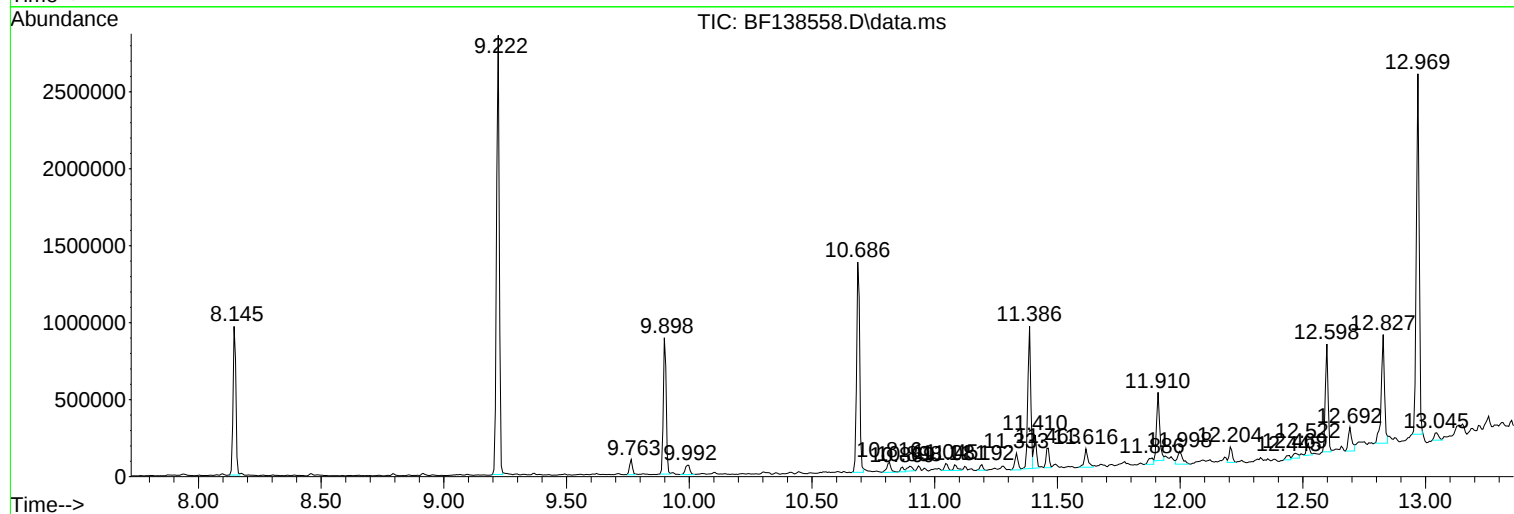
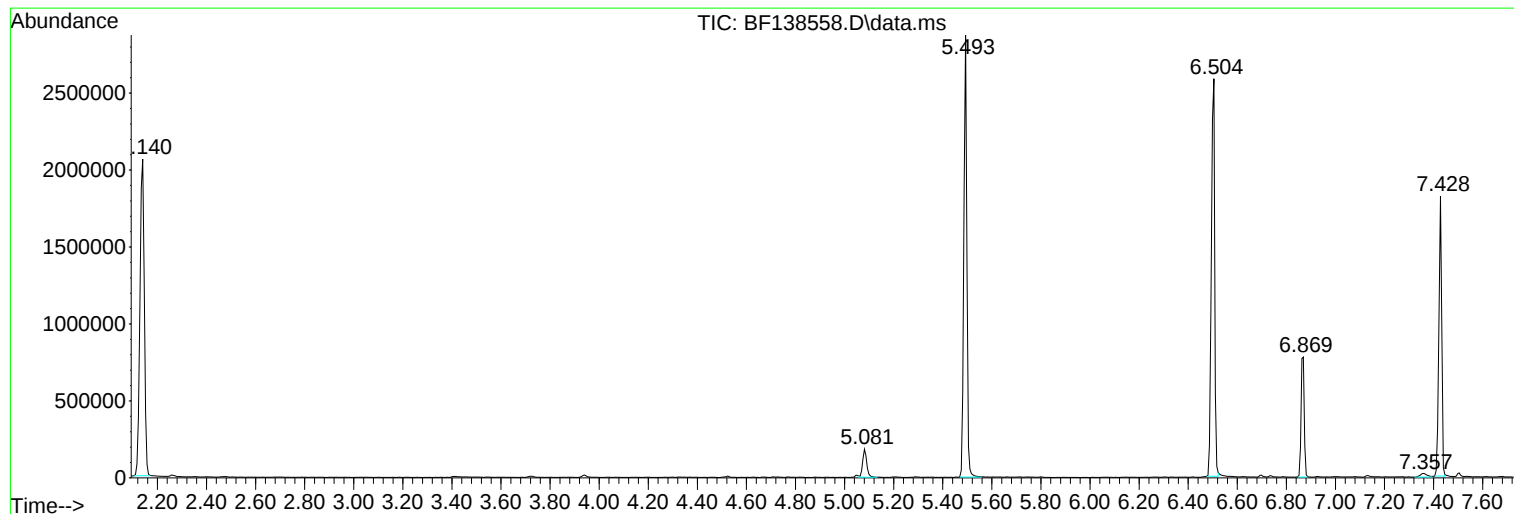
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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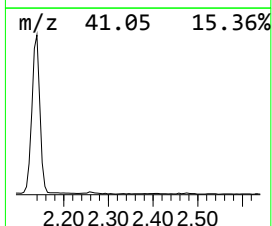
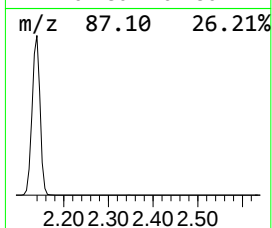
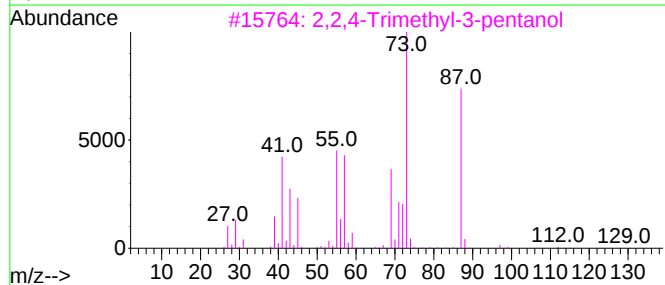
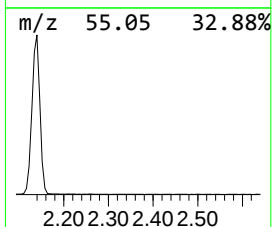
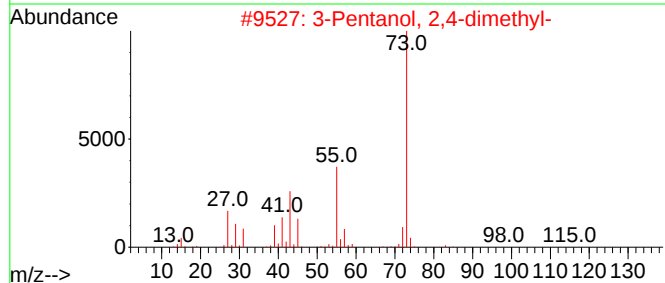
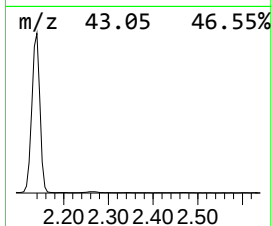
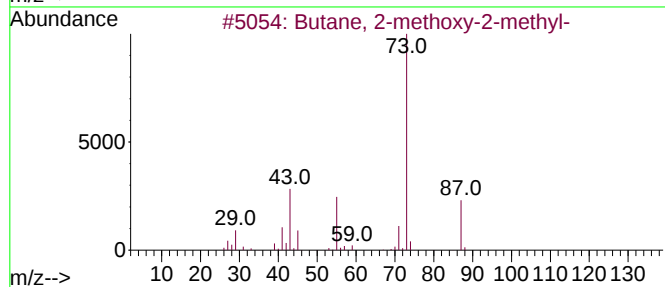
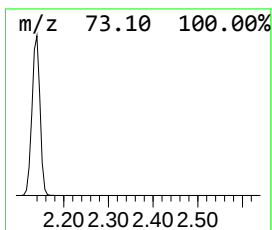
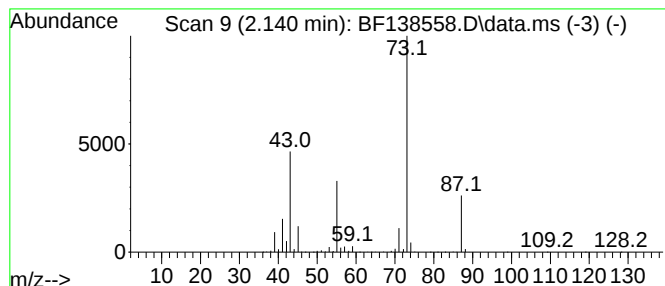
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Butane, 2-methoxy-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.140	73.25 ng	2557380	1,4-Dichlorobenzene-d4	6.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			3-Pentanol, 2,4-dimethyl-	116	C7H16O	000600-36-2	43
3			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
4			Silane, tetramethyl-	88	C4H12Si	000075-76-3	38
5			1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	25



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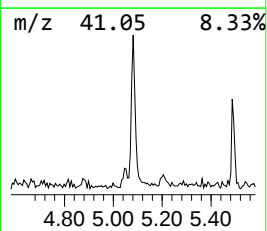
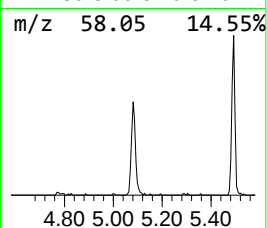
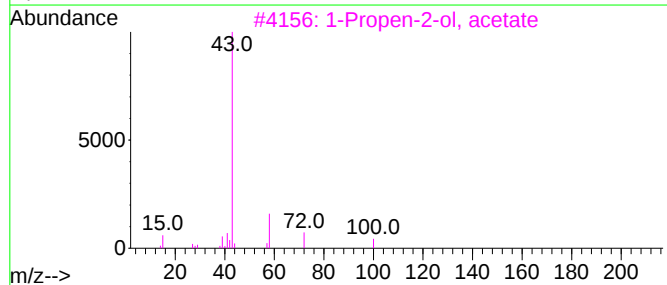
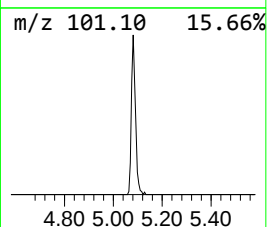
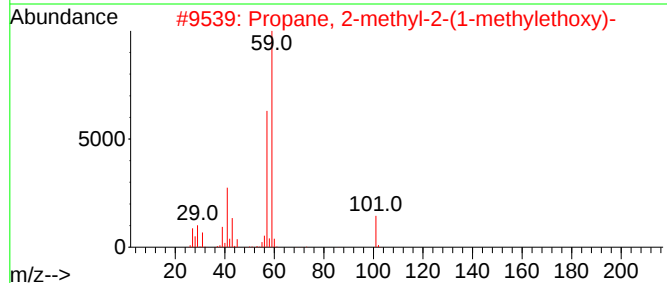
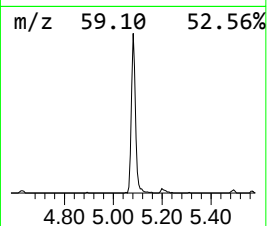
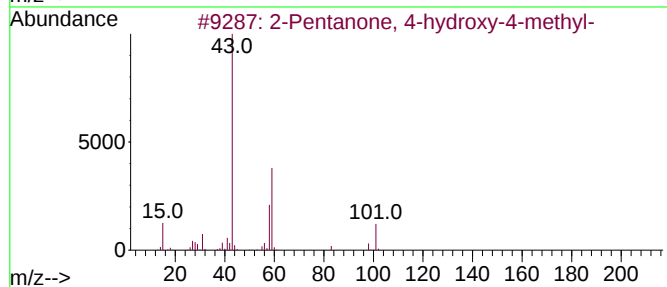
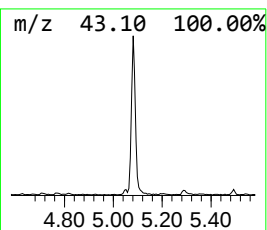
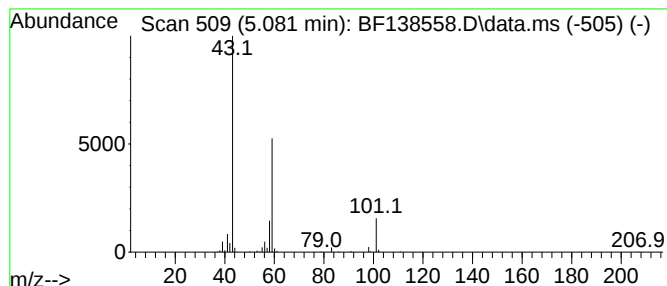
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.081	6.00 ng	209664	1,4-Dichlorobenzene-d4	6.869

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	59
2		Propane, 2-methyl-2-(1-methyleth...	116	C7H16O	017348-59-3	36
3		1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
4		2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	9
5		3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	9



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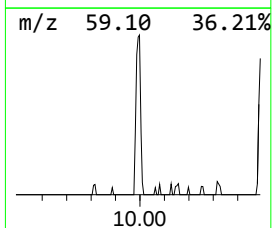
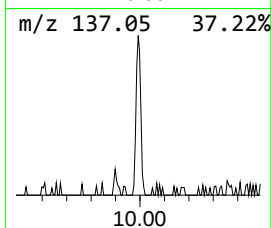
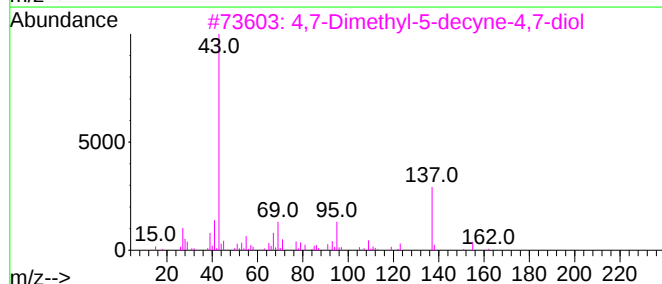
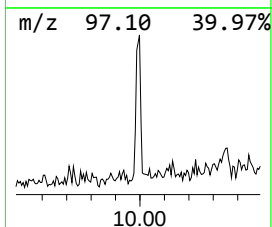
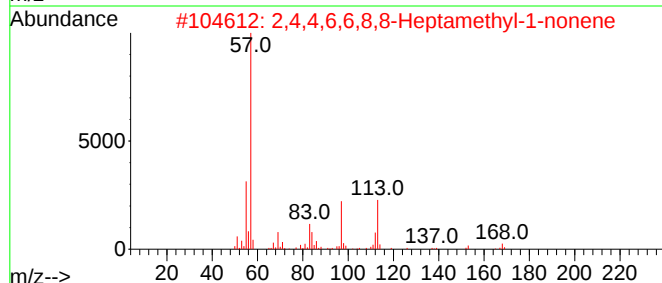
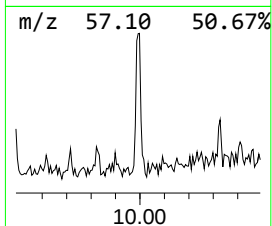
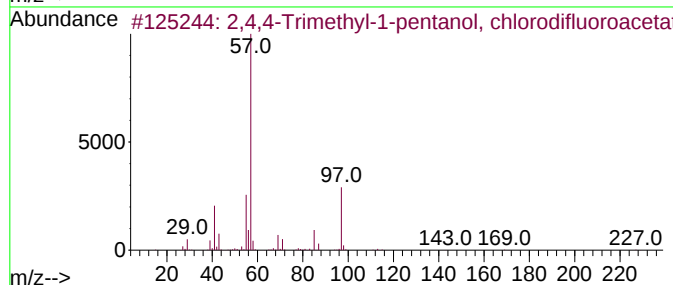
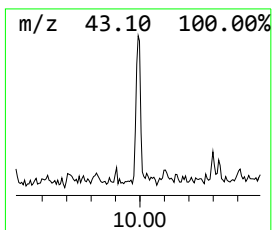
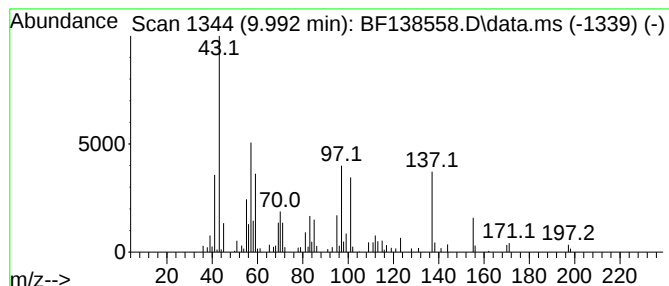
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown9.992 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.992	2.22 ng	80599	Acenaphthene-d10	9.898

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,4,4-Trimethyl-1-pentanol, chlo...	242	C10H17ClF2O2	1000447-27-2	12		
2	2,4,4,6,6,8,8-Heptamethyl-1-nonene	224	C16H32	015796-04-0	10		
3	4,7-Dimethyl-5-decyne-4,7-diol	198	C12H22O2	000126-87-4	10		
4	Carbonic acid, heptadecyl prop-1...	340	C21H40O3	1000382-90-2	10		
5	6,10,13-Trimethyltetradecanol	256	C17H36O	1000131-71-0	10		



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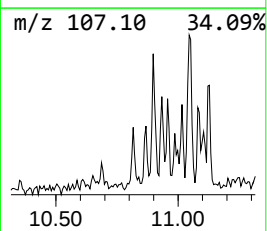
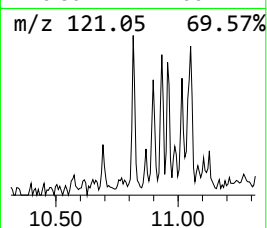
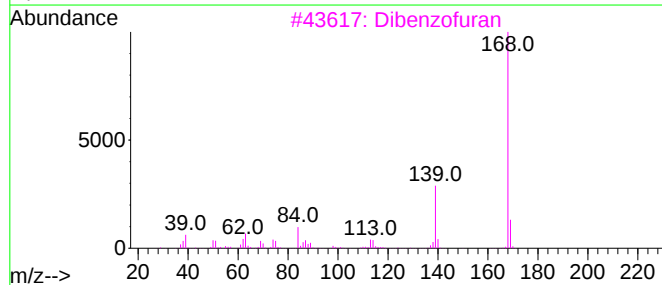
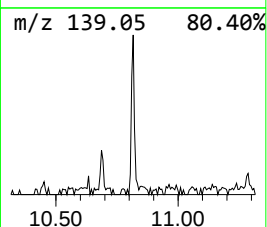
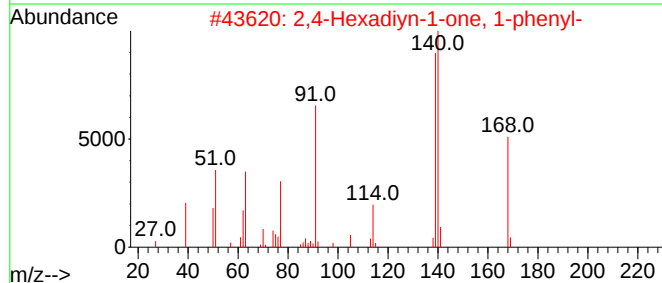
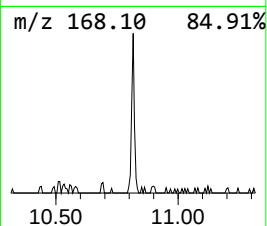
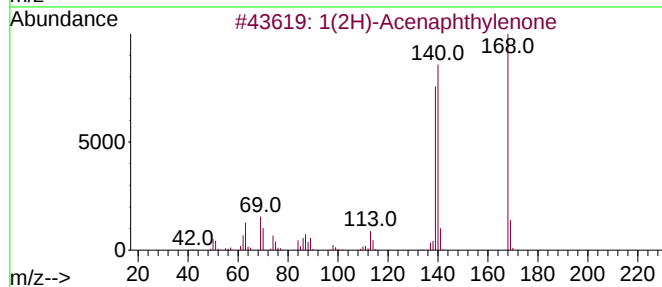
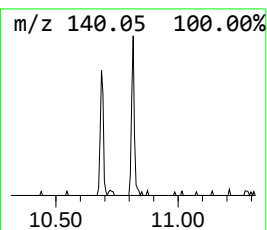
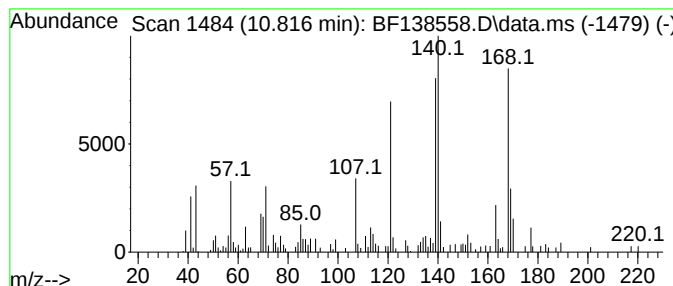
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 1(2H)-Acenaphthylenone Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	2.08 ng	79047	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	93
2			2,4-Hexadiyn-1-one, 1-phenyl-	168	C12H8O	000495-74-9	64
3			Dibenzofuran	168	C12H8O	000132-64-9	43
4			Ethyl maltol	140	C7H8O3	004940-11-8	43
5			Cyclobuta[de]naphthalene	140	C11H8	024973-91-9	43



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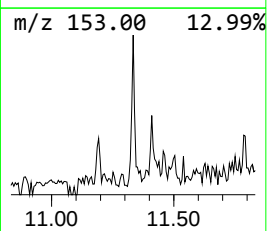
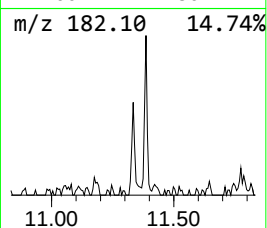
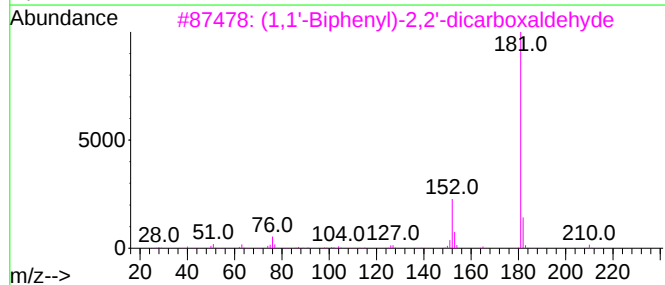
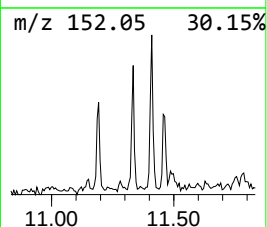
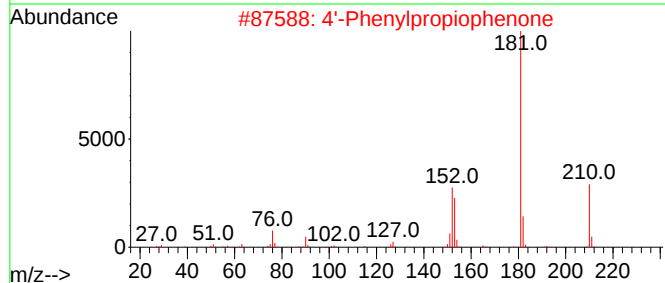
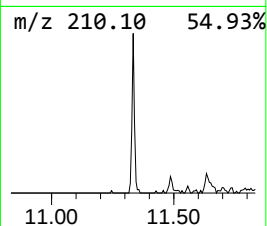
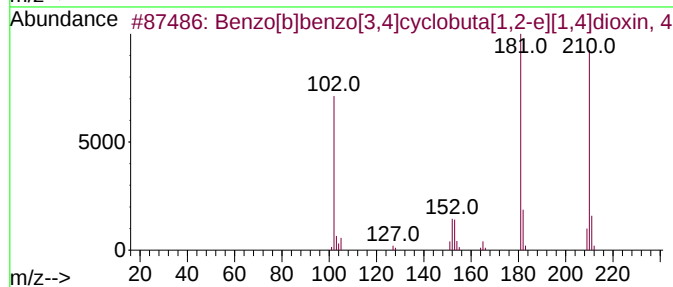
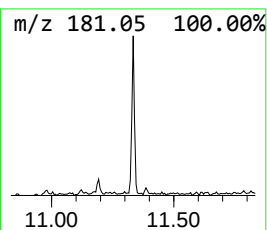
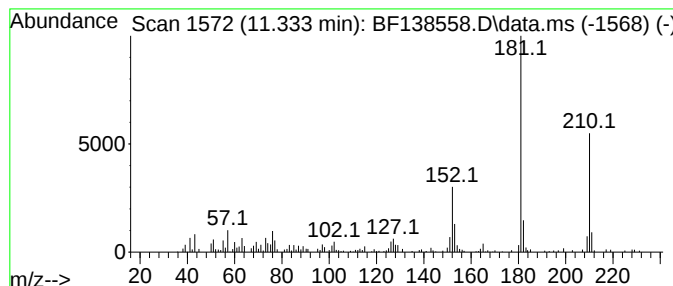
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Benzo[b]benzo[3,4]cyclobuta... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.334	2.65 ng	100528	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]benzo[3,4]cyclobuta[1,2-...	210	C14H10O2	042896-18-4	93
2			4'-Phenylpropiophenone	210	C15H14O	037940-57-1	87
3			(1,1'-Biphenyl)-2,2'-dicarboxald...	210	C14H10O2	001210-05-5	50
4			2-Fluorenamine	181	C13H11N	000153-78-6	49
5			3-Methyl-2-azafluorene	181	C13H11N	018631-12-4	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-3

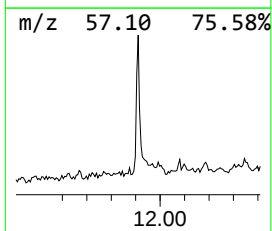
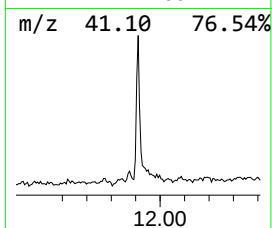
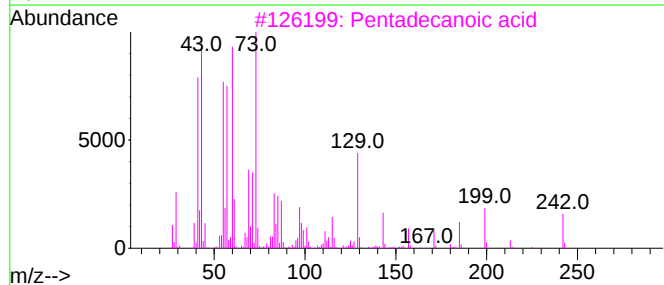
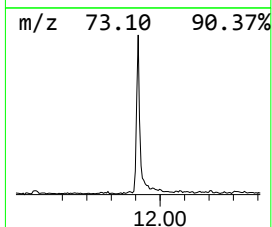
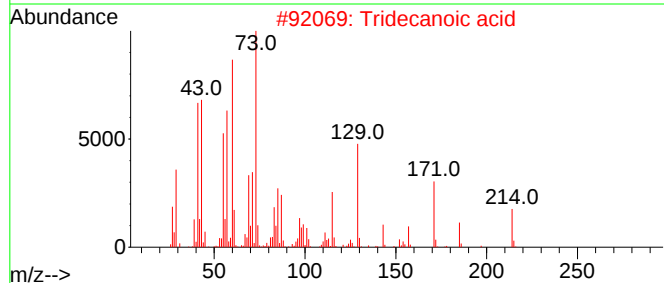
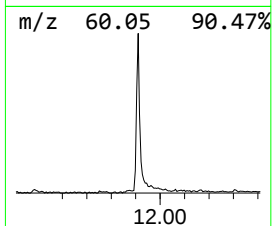
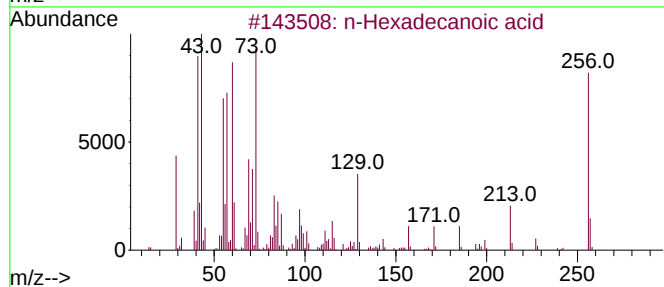
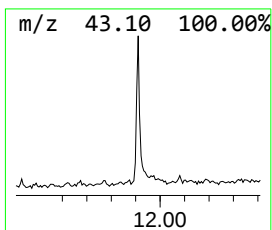
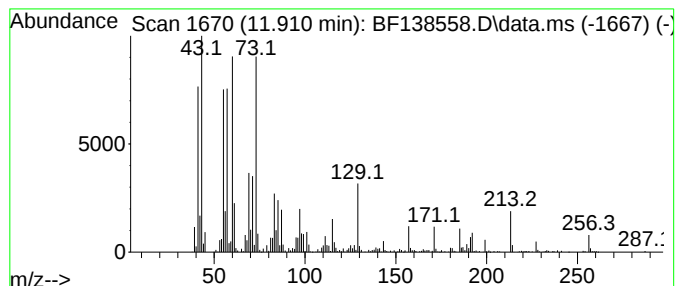
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.910	10.58 ng	401776	Phenanthrene-d10	11.386

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	96
3			Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4			Octadecanoic acid	284	C18H36O2	000057-11-4	70
5			8-Methylnonanoic acid	172	C10H20O2	005963-14-4	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

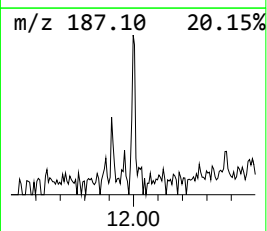
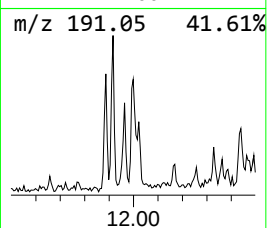
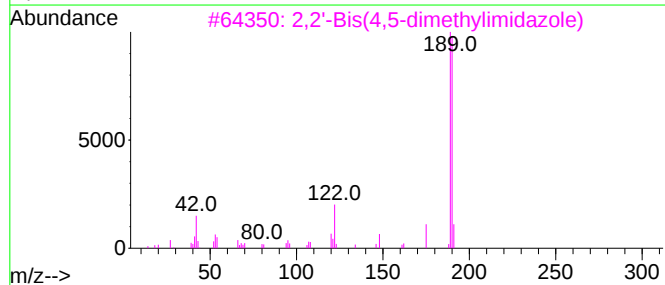
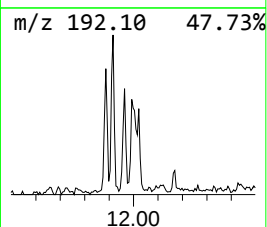
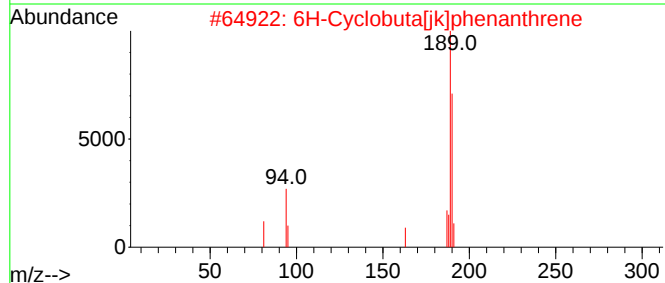
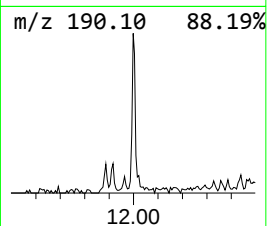
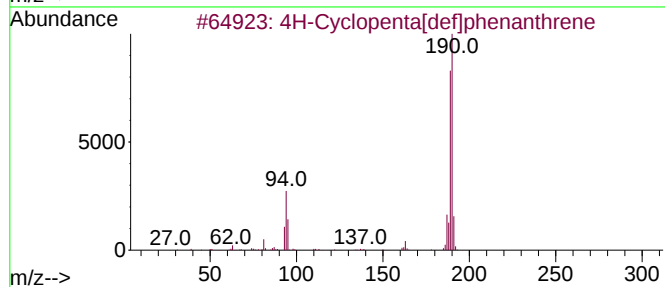
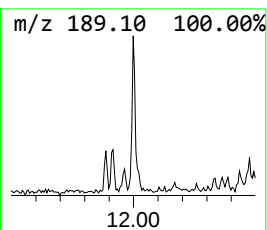
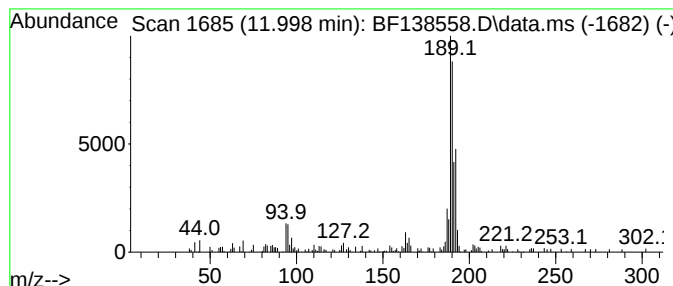
Instrument :
BNA_F
ClientSampleId :
WC-3

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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 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.998	3.09 ng	117439	Phenanthrene-d10	11.386	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	68
2	6H-Cyclobuta[jk]phenanthrene	190	C15H10	083469-43-6	58
3	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	43
4	Carbonic acid, ethyl 2-formyl-4,...	262	C10H8Cl2O4	1000331-34-4	37
5	6-Amino-2,4-dichloro-3-ethylphenol	205	C8H9Cl2NO	101819-99-2	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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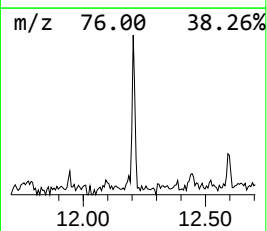
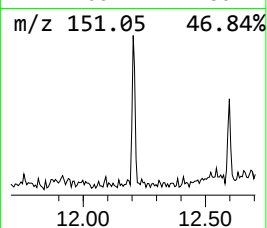
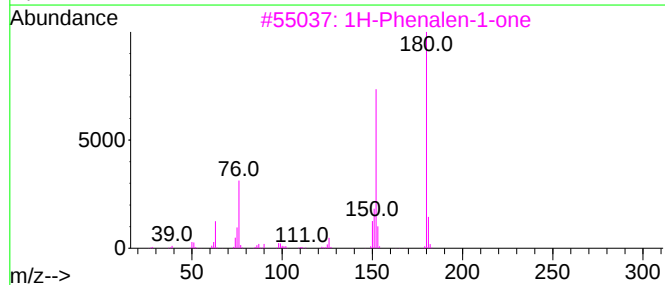
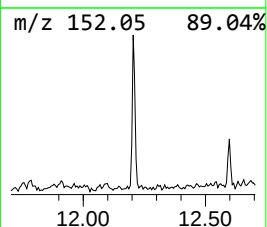
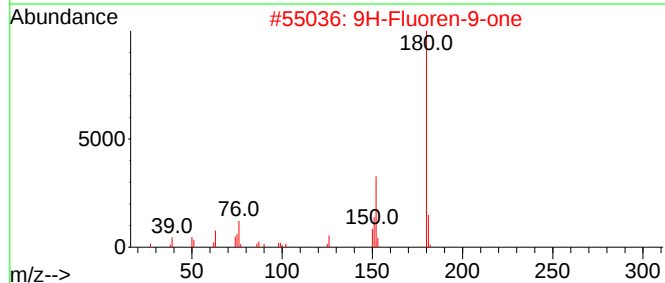
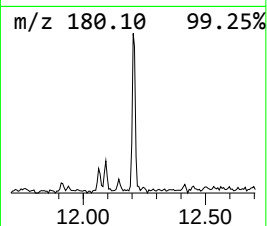
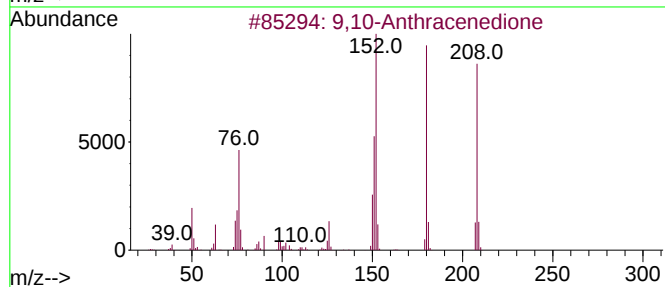
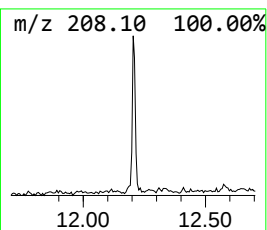
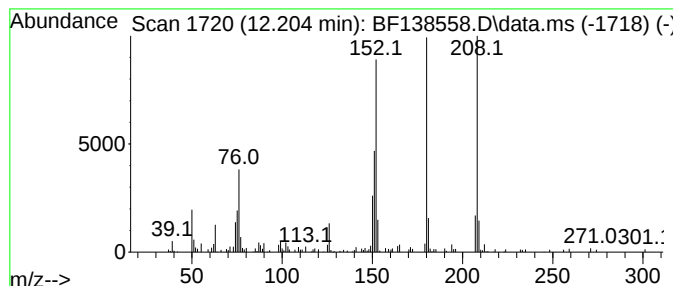
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 9,10-Anthracenedione Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.204	2.28 ng	86694	Phenanthrene-d10	11.386

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9,10-Anthracenedione	208	C14H8O2	000084-65-1	96
2		9H-Fluoren-9-one	180	C13H8O	000486-25-9	78
3		1H-Phenalen-1-one	180	C13H8O	000548-39-0	78
4		Benzo[c]cinnoline	180	C12H8N2	000230-17-1	78
5		Benzo[h]cinnoline	180	C12H8N2	000230-31-9	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-3

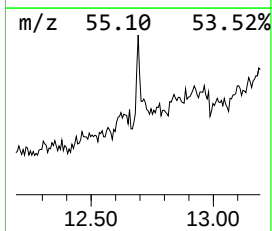
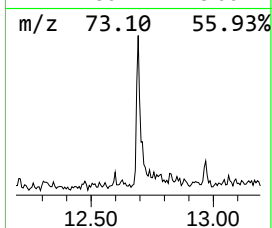
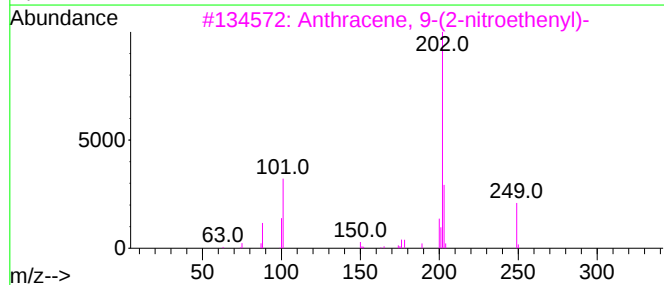
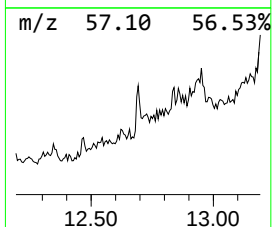
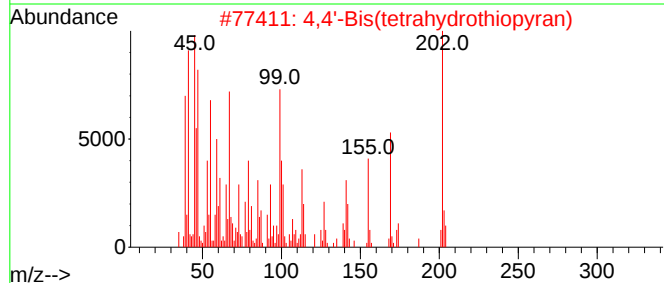
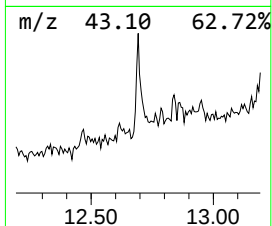
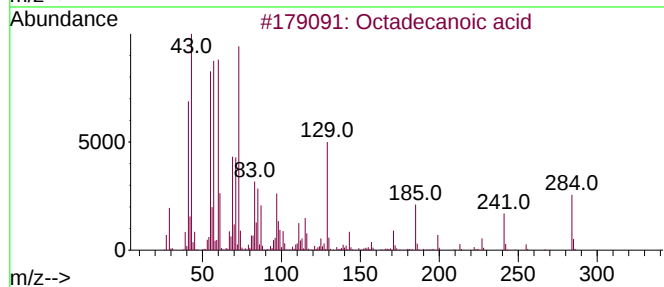
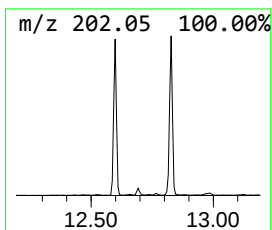
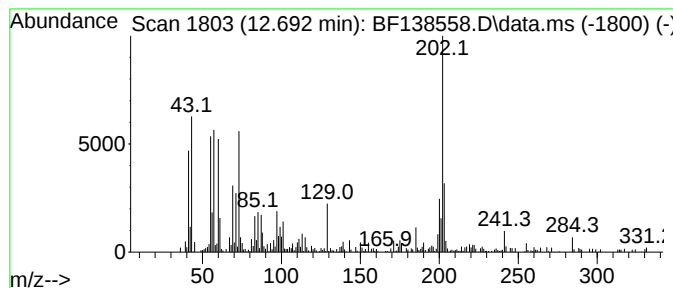
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.692	4.18 ng	158772	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	93
2			4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	92
3			Anthracene, 9-(2-nitroethenyl)-	249	C16H11NO2	058349-77-2	53
4			Dodecanoic acid	200	C12H24O2	000143-07-7	49
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	48



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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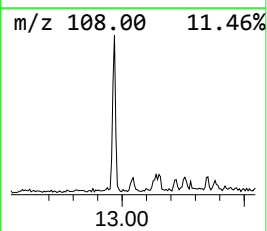
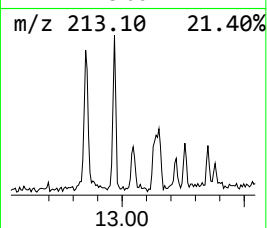
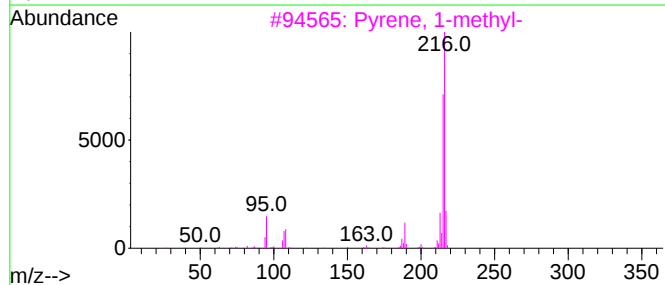
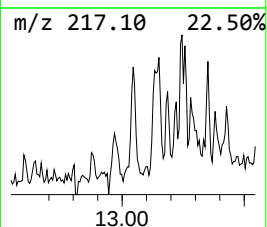
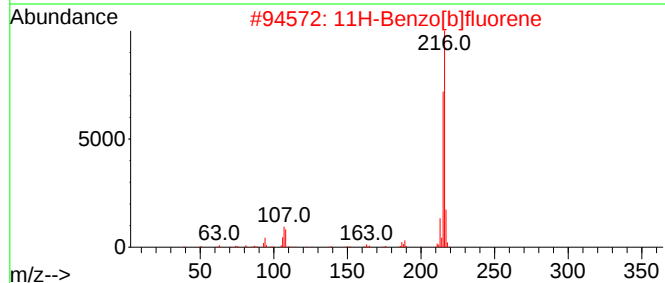
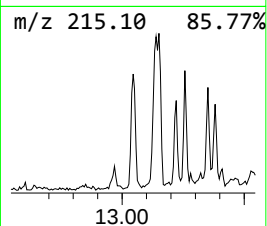
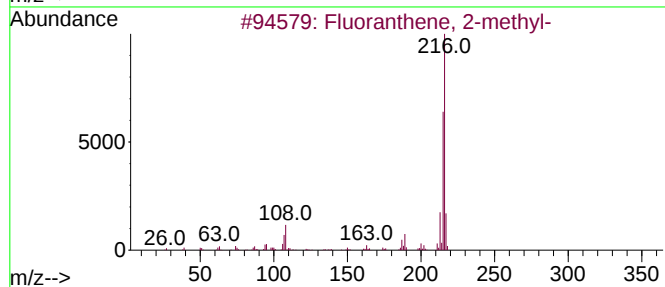
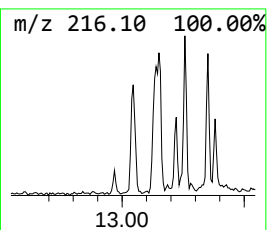
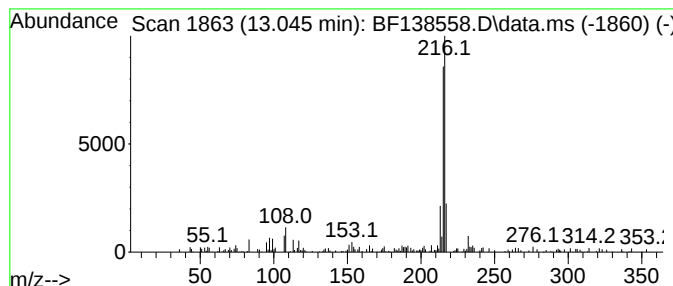
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Fluoranthene, 2-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.045	2.03 ng	62060	Chrysene-d12	14.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	91
2			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	87
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	72
4			3-Pyridinamine, 6-chloro-N-(phen...	216	C12H9ClN2	005325-71-3	64
5			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
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Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

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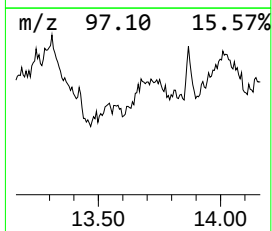
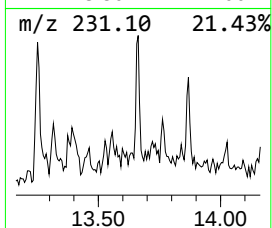
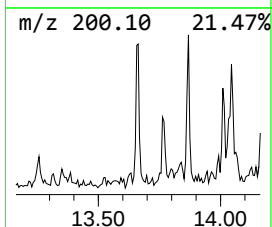
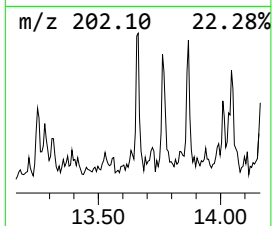
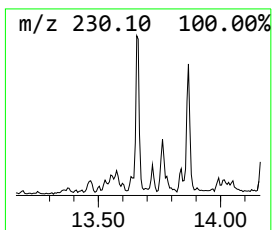
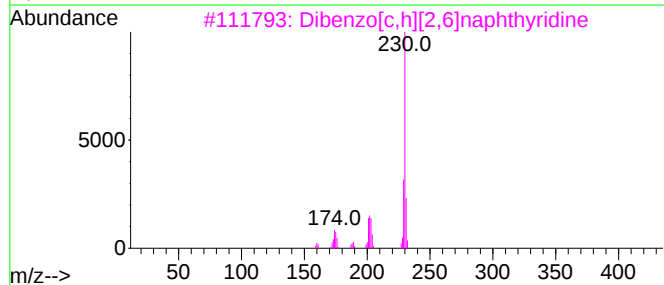
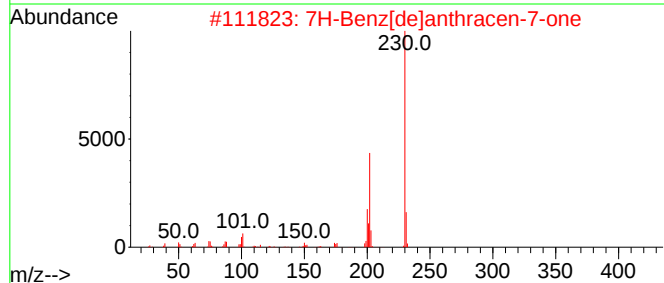
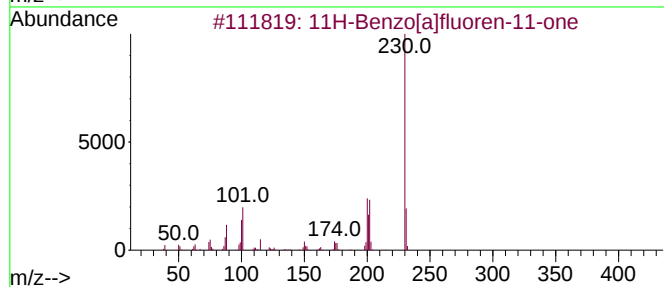
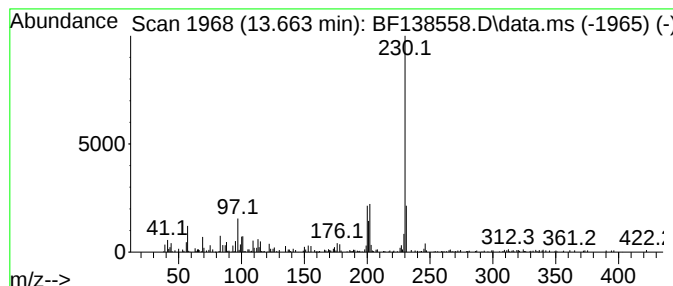
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 11H-Benzo[a]fluoren-11-one Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	4.27 ng	130316	Chrysene-d12	14.022

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	83
2			7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	76
3			Dibenzo[c,h][2,6]naphthyridine	230	C16H10N2	000218-30-4	59
4			(E)-.alpha.,.alpha.'-Dicyanostil...	230	C16H10N2	002450-55-7	53
5			benzonitrile, 2-(3-isoquinolinyl)-	230	C16H10N2	1000401-00-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

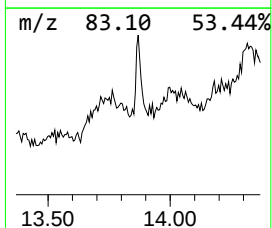
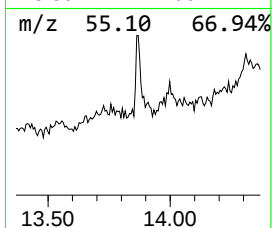
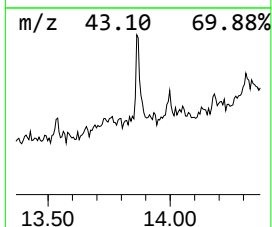
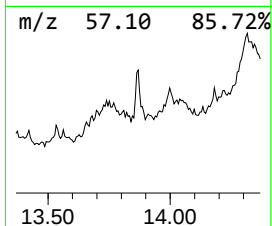
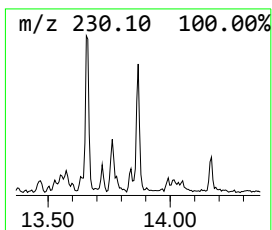
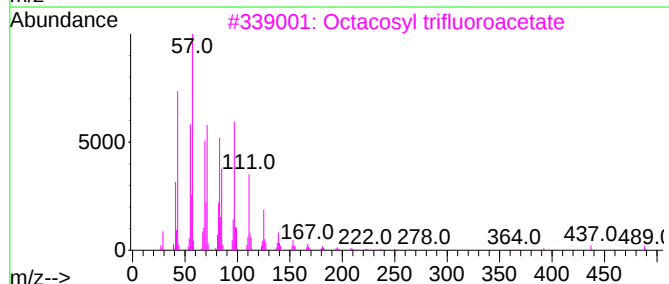
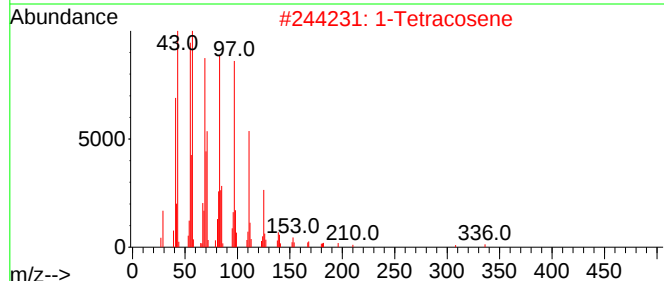
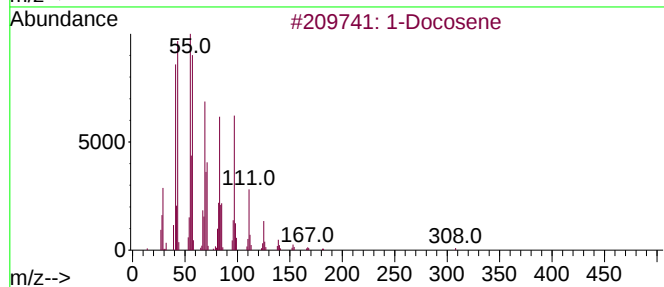
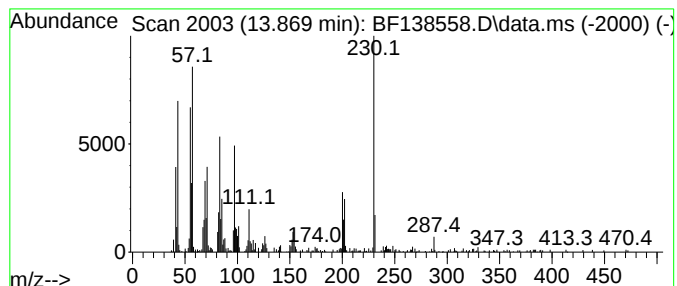
Instrument :
BNA_F
ClientSampleId :
WC-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 1-Docosene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.869	6.67 ng	203640	Chrysene-d12	14.022	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	89
2	1-Tetracosene	336	C24H48	010192-32-2	64
3	Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	50
4	Cyclotetracosane	336	C24H48	000297-03-0	46
5	Carbonic acid, hexadecyl isobuty...	342	C21H42O3	1010314-61-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-3

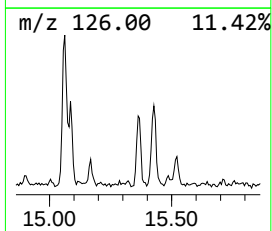
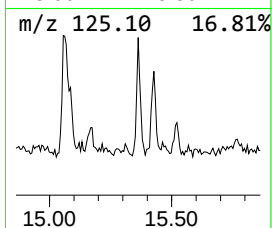
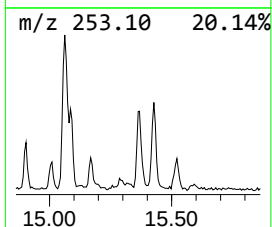
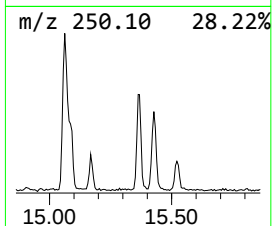
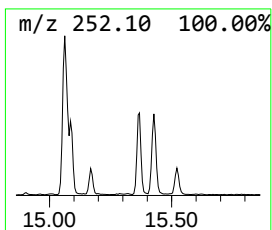
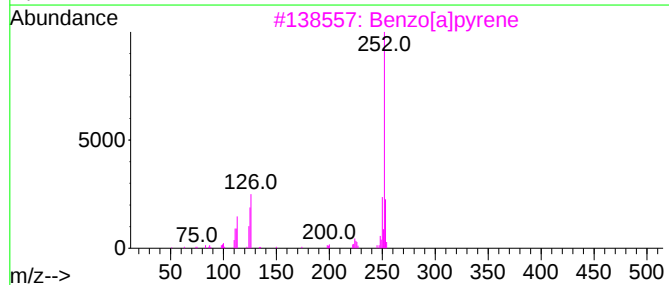
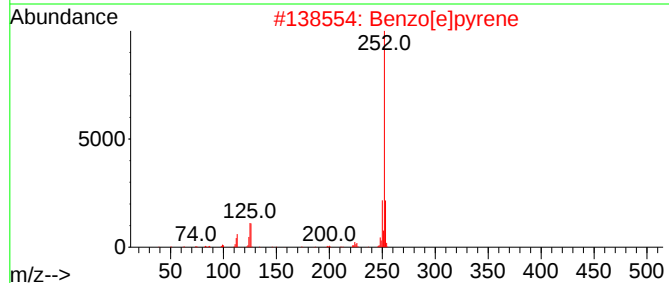
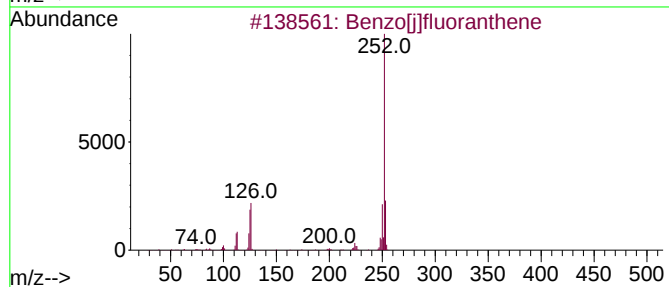
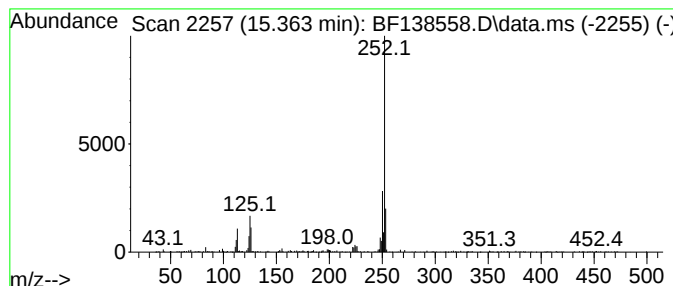
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Benzo[j]fluoranthene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.363	11.20 ng	209329	Perylene-d12	15.492

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzo[j]fluoranthene	252	C20H12	000205-82-3	93
2		Benzo[e]pyrene	252	C20H12	000192-97-2	93
3		Benzo[a]pyrene	252	C20H12	000050-32-8	91
4		Benzo[k]fluoranthene	252	C20H12	000207-08-9	91
5		Benzo[b]fluoranthene	252	C20H12	000205-99-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF071224\
Data File : BF138558.D
Acq On : 12 Jul 2024 21:09
Operator : CG\JU
Sample : P3187-09
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF070924.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
Butane, 2-metho...	2.140	73.3	ng	2557380	1	6.869	698299	20.0
2-Pentanone, 4-...	5.081	6.0	ng	209664	1	6.869	698299	20.0
unknown9.992	9.992	2.2	ng	80599	3	9.898	727173	20.0
1(2H)-Acenaphth...	10.816	2.1	ng	79047	4	11.386	759631	20.0
Benzo[b]benzo[3...	11.334	2.6	ng	100528	4	11.386	759631	20.0
n-Hexadecanoic ...	11.910	10.6	ng	401776	4	11.386	759631	20.0
4H-Cyclopenta[d...	11.998	3.1	ng	117439	4	11.386	759631	20.0
9,10-Anthracene...	12.204	2.3	ng	86694	4	11.386	759631	20.0
Octadecanoic acid	12.692	4.2	ng	158772	4	11.386	759631	20.0
Fluoranthene, 2...	13.045	2.0	ng	62060	5	14.022	610491	20.0
11H-Benzo[a]flu...	13.663	4.3	ng	130316	5	14.022	610491	20.0
1-Docosene	13.869	6.7	ng	203640	5	14.022	610491	20.0
Benzo[j]fluoran...	15.363	11.2	ng	209329	6	15.492	373742	20.0