

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
 Data File : BF139197.D
 Acq On : 27 Aug 2024 20:28
 Operator : RC/JU
 Sample : P3732-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 UI-1

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-BF082724.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF139197.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.181	8	15	22	rVB	1046894	1647122	100.00%	8.518%
2	5.104	503	512	521	rBV	67168	106265	6.45%	0.550%
3	5.504	574	580	585	rBV	1204694	1584876	96.22%	8.196%
4	6.510	745	751	757	rBV	1218412	1635532	99.30%	8.458%
5	6.710	781	785	793	rBV	22581	31186	1.89%	0.161%
6	6.892	811	816	821	rBV	452869	547318	33.23%	2.831%
7	7.451	905	911	916	rBV	802747	993492	60.32%	5.138%
8	7.704	950	954	958	rBV	15904	18159	1.10%	0.094%
9	8.175	1028	1034	1041	rBV	532186	674206	40.93%	3.487%
10	8.481	1081	1086	1095	rBV	61345	77214	4.69%	0.399%
11	9.245	1211	1216	1221	rBV	1366364	1645048	99.87%	8.508%
12	9.786	1304	1308	1313	rBV	36763	45549	2.77%	0.236%
13	9.922	1326	1331	1336	rBV	550989	675702	41.02%	3.494%
14	10.022	1341	1348	1361	rBV	606316	931066	56.53%	4.815%
15	10.710	1459	1465	1476	rVB	623780	790189	47.97%	4.087%
16	10.833	1481	1486	1493	rVB4	16195	25223	1.53%	0.130%
17	11.169	1535	1543	1547	rBV6	7028	17896	1.09%	0.093%
18	11.410	1579	1584	1592	rBV2	437638	709439	43.07%	3.669%
19	11.480	1592	1596	1600	rBV	28739	35130	2.13%	0.182%
20	11.798	1645	1650	1657	rVB6	9555	18037	1.10%	0.093%
21	11.910	1665	1669	1670	rBV	34972	39992	2.43%	0.207%
22	11.933	1670	1673	1678	rVB2	72054	96913	5.88%	0.501%
23	12.022	1684	1688	1697	rVB	59461	115882	7.04%	0.599%
24	12.204	1715	1719	1729	rVB3	24616	56487	3.43%	0.292%
25	12.357	1738	1745	1747	rBV2	11823	21004	1.28%	0.109%
26	12.457	1757	1762	1765	rBV	45789	62596	3.80%	0.324%
27	12.498	1765	1769	1773	rVV4	29500	60313	3.66%	0.312%
28	12.539	1773	1776	1783	rVB2	35593	55114	3.35%	0.285%
29	12.621	1785	1790	1794	rVV	429018	528416	32.08%	2.733%
30	12.663	1794	1797	1803	rBV5	11987	26923	1.63%	0.139%
31	12.716	1803	1806	1809	rVV	20943	23643	1.44%	0.122%
32	12.774	1813	1816	1822	rVV3	11387	18719	1.14%	0.097%
33	12.851	1822	1829	1834	rVV	426694	614717	37.32%	3.179%
34	12.898	1834	1837	1842	rVB2	23162	34854	2.12%	0.180%
35	12.992	1845	1853	1860	rVB	1001788	1301280	79.00%	6.730%
36	13.069	1860	1866	1873	rBV3	52274	109228	6.63%	0.565%

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37	13.174	1873	1884	1888	rBV2	66427	143759	8.73%	0.743%
38	13.239	1888	1895	1898	rBV3	39681	67910	4.12%	0.351%
39	13.280	1898	1902	1906	rVV2	52224	63900	3.88%	0.330%
40	13.374	1914	1918	1920	rBV	32882	38743	2.35%	0.200%
41	13.404	1920	1923	1927	rVB	30527	35097	2.13%	0.182%
42	13.486	1933	1937	1941	rBV6	13013	24500	1.49%	0.127%
43	13.680	1966	1970	1976	rVB	45402	67053	4.07%	0.347%
44	13.792	1984	1989	1992	rBV2	54040	99159	6.02%	0.513%
45	13.827	1992	1995	1996	rBV	25165	28822	1.75%	0.149%
46	13.892	2002	2006	2013	rVB2	89974	137607	8.35%	0.712%
47	14.045	2025	2032	2035	rBV2	567456	990258	60.12%	5.121%
48	14.151	2047	2050	2053	rVB	29212	32301	1.96%	0.167%
49	14.186	2053	2056	2059	rBV	20331	28599	1.74%	0.148%
50	14.427	2094	2097	2101	rVB	42296	53359	3.24%	0.276%
51	14.463	2101	2103	2107	rVB	33417	38380	2.33%	0.198%
52	14.521	2110	2113	2116	rVV2	38822	55475	3.37%	0.287%
53	14.557	2116	2119	2127	rVB2	42188	92253	5.60%	0.477%
54	15.092	2204	2210	2220	rVB	271833	619900	37.64%	3.206%
55	15.198	2225	2228	2232	rVB	48305	65370	3.97%	0.338%
56	15.392	2257	2261	2267	rVB	135331	206640	12.55%	1.069%
57	15.457	2267	2272	2277	rVB	200986	291830	17.72%	1.509%
58	15.521	2277	2283	2287	rBV	219560	363228	22.05%	1.878%
59	16.992	2527	2533	2543	rVB2	98341	214479	13.02%	1.109%
60	17.439	2603	2609	2621	rVB	75461	172572	10.48%	0.892%
61	17.668	2645	2648	2656	rVB	15601	30404	1.85%	0.157%

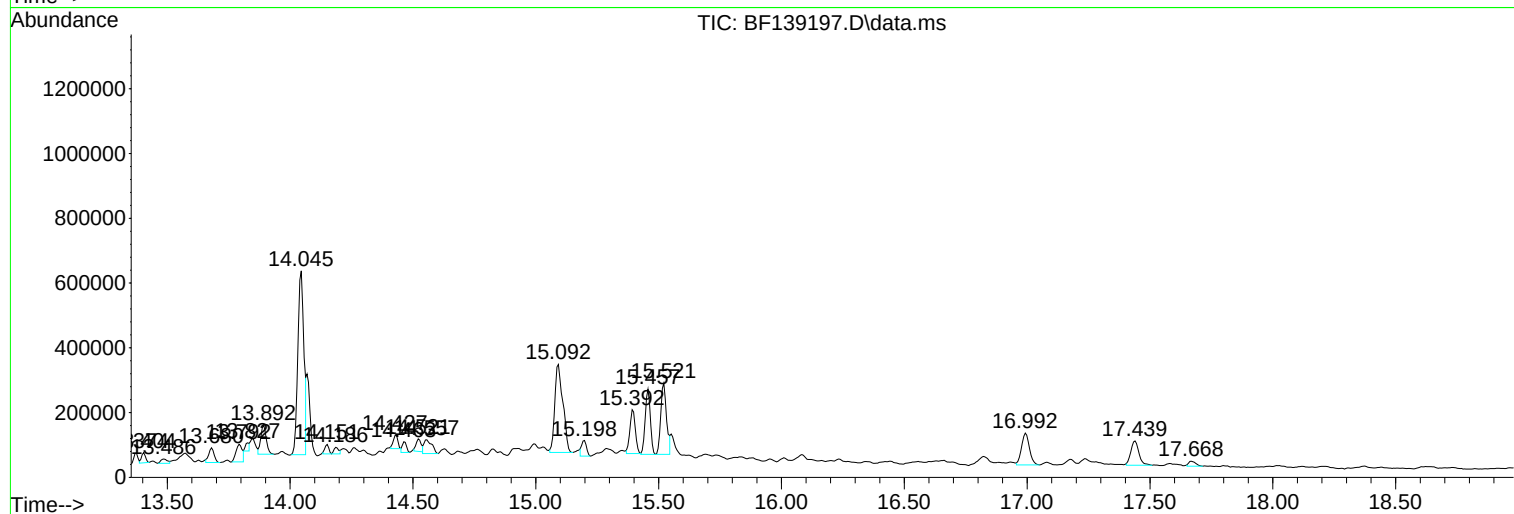
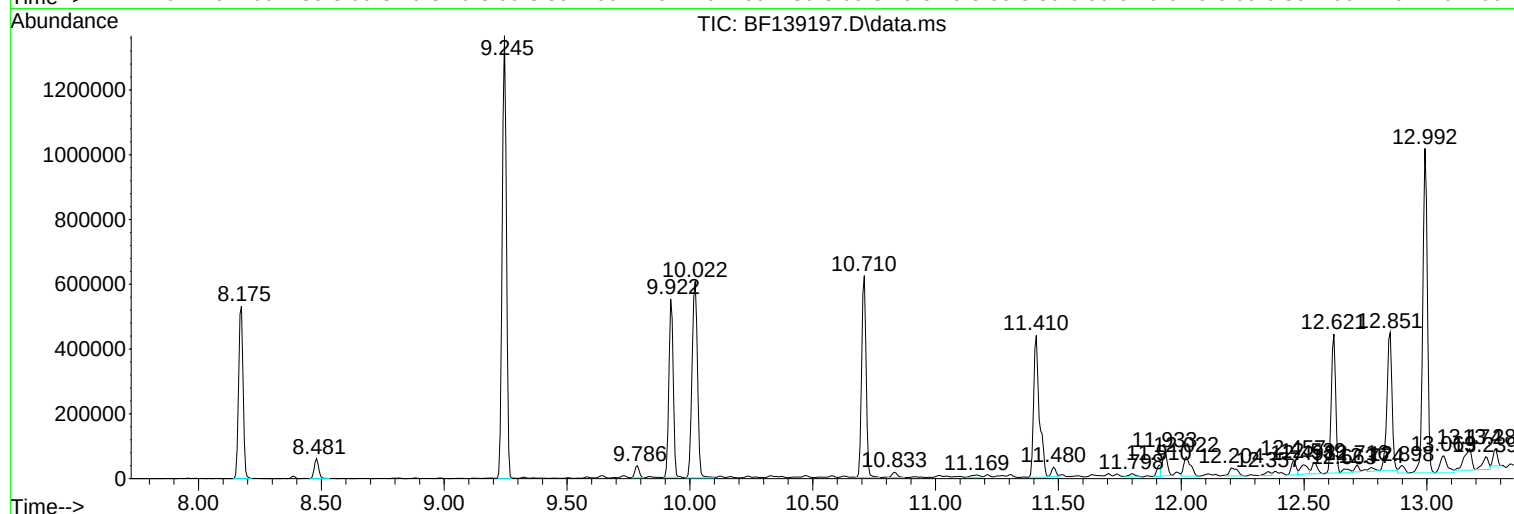
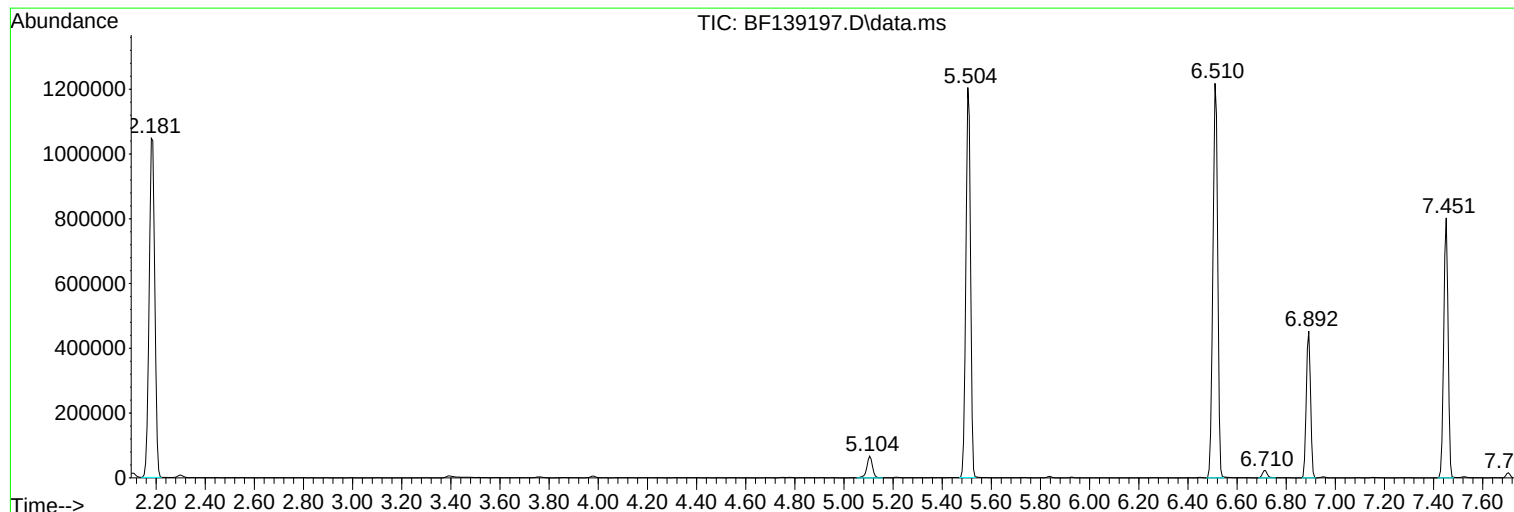
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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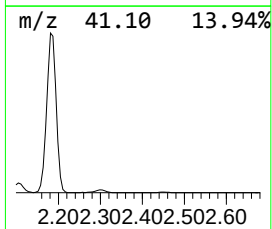
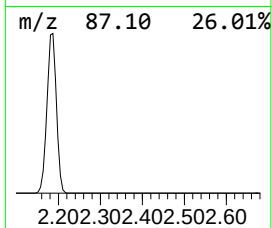
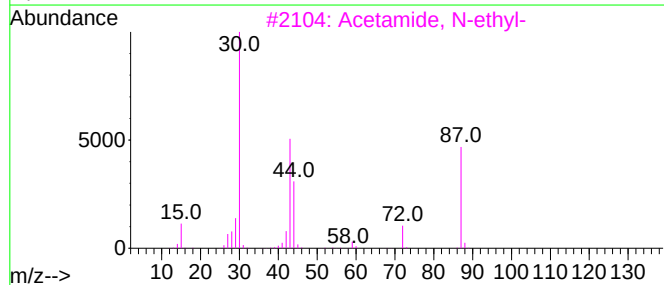
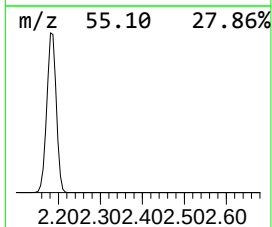
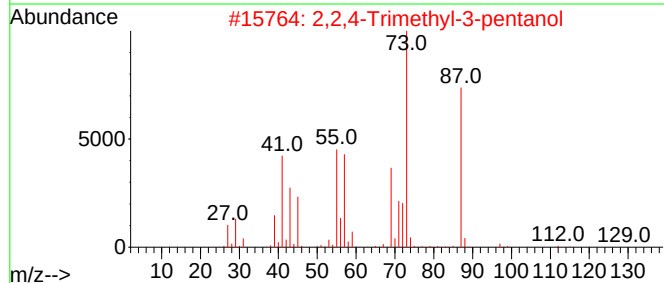
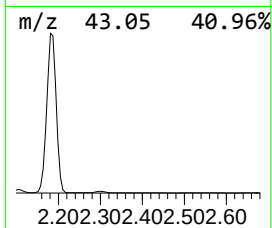
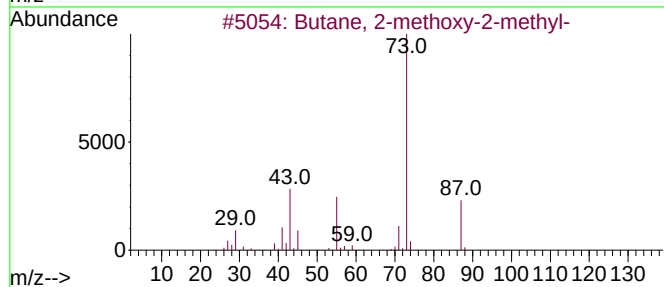
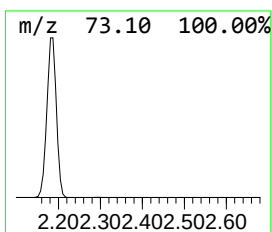
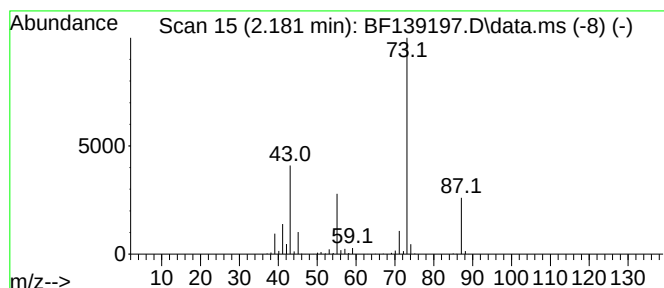
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082724.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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.
2.181	60.19 ng	1647120	1,4-Dichlorobenzene-d4	6.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	27
4			Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
5			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12



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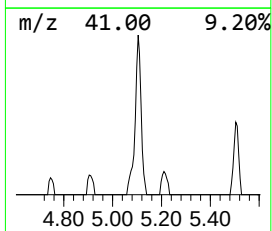
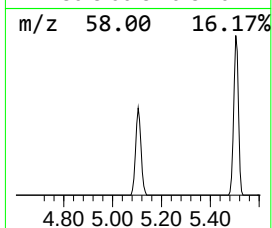
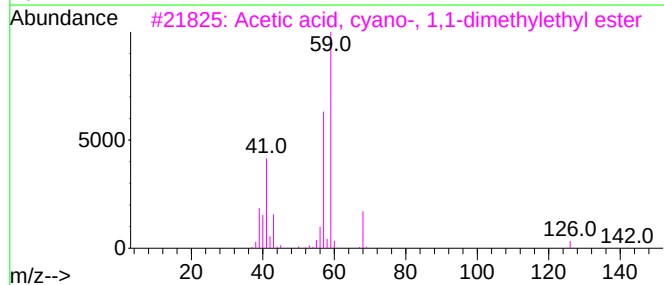
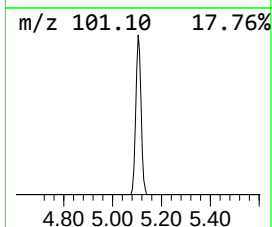
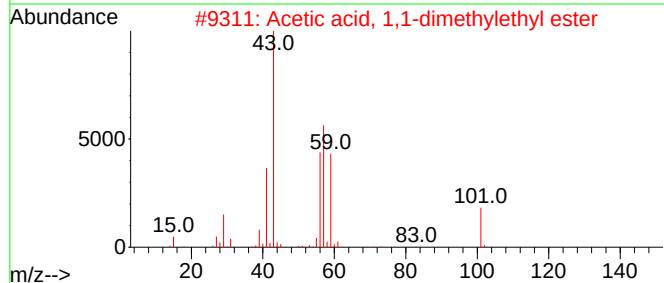
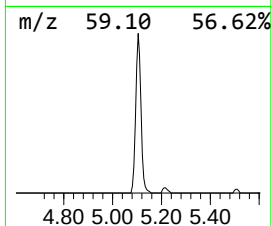
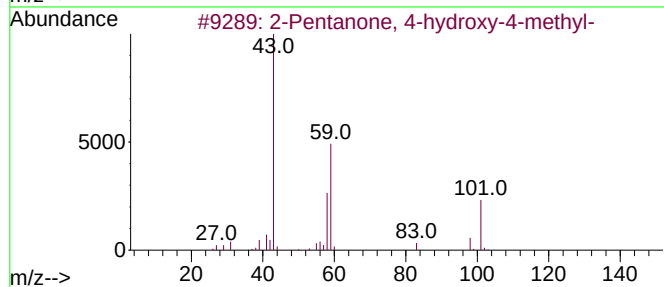
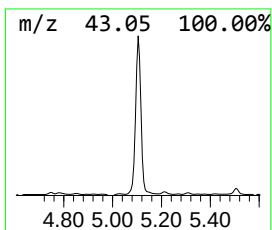
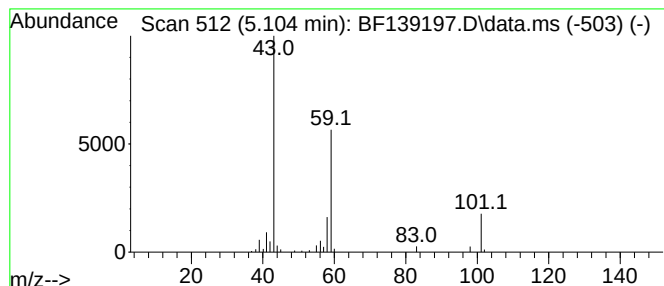
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.104	3.88 ng	106265	1,4-Dichlorobenzene-d4	6.892

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56
2			Acetic acid, 1,1-dimethylethyl e...	116	C6H12O2	000540-88-5	38
3			Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	25
4			1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12
5			2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	10



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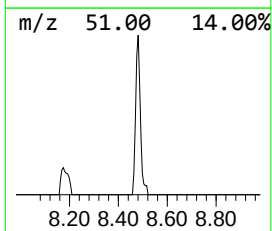
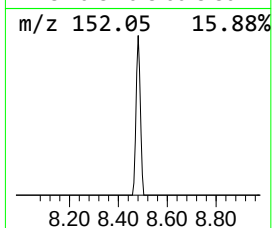
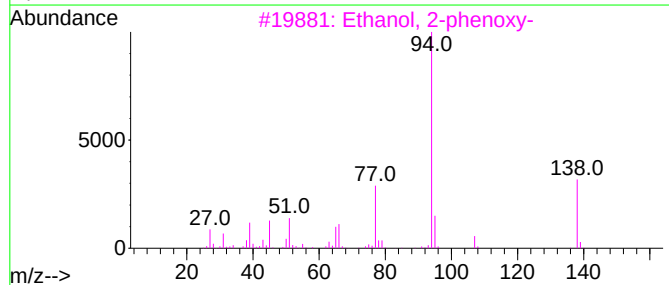
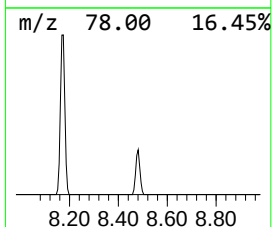
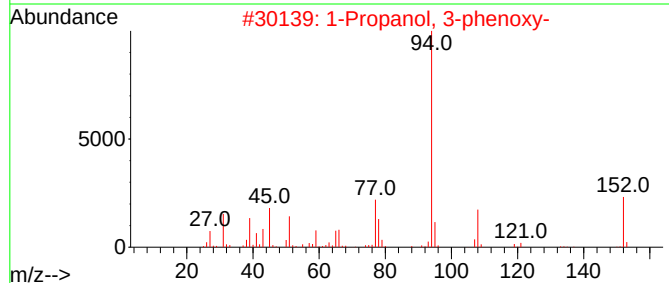
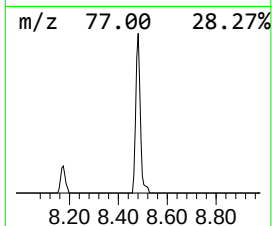
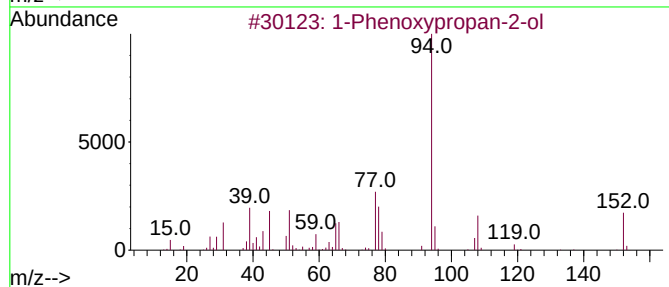
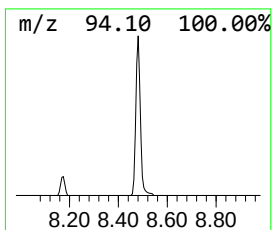
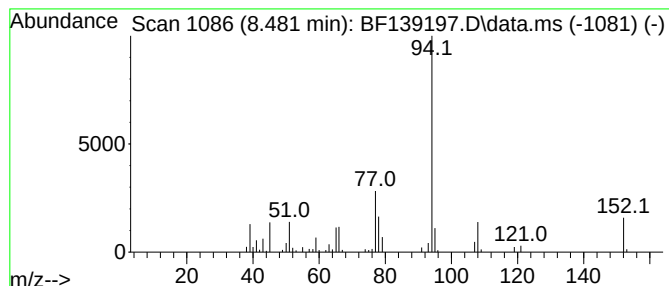
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Phenoxypropan-2-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.481	2.29 ng	77214	Naphthalene-d8	8.175

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	87
3			Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4			Carbonic acid, neopentyl phenyl ...	208	C12H16O3	013183-19-2	59
5			1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59



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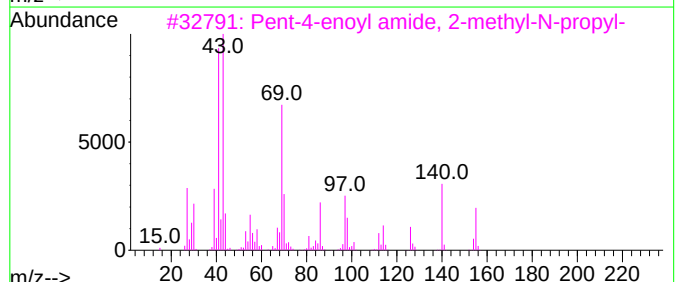
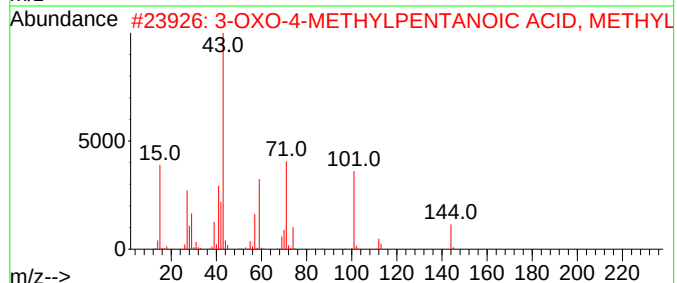
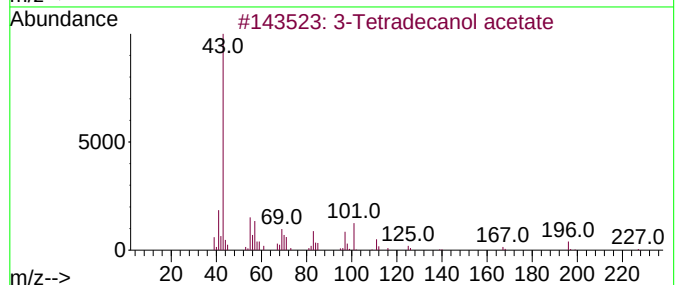
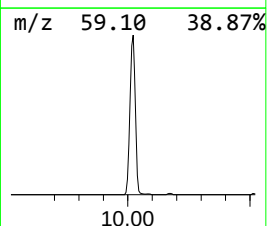
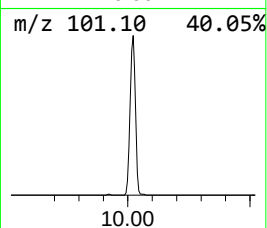
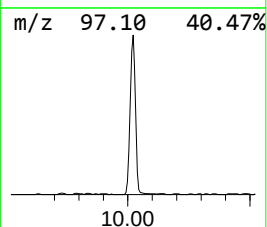
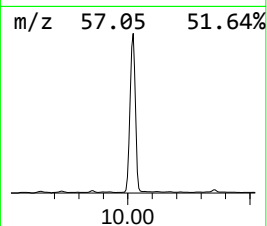
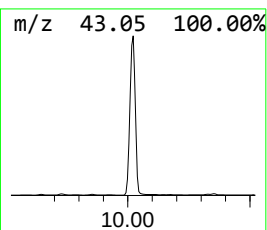
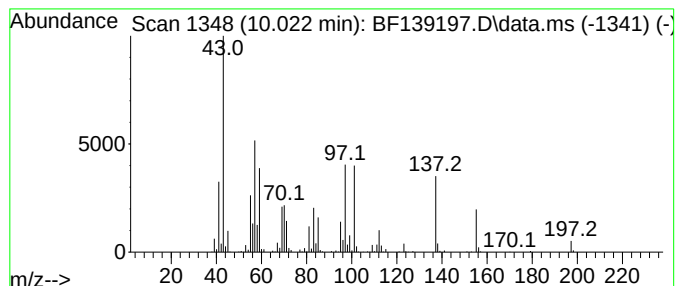
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 unknown10.022 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.022	27.56 ng	931066	Acenaphthene-d10	9.922

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Tetradecanol acetate	256	C16H32O2	051354-23-5	25
2			3-OXO-4-METHYLPENTANOIC ACID, ME...	144	C7H12O3	042558-54-3	16
3			Pent-4-enoyl amide, 2-methyl-N-p...	155	C9H17NO	1000420-35-2	14
4			11-Tricosene	322	C23H46	052078-56-5	12
5			3-Octanol, 2,6-dimethyl-, acetate	200	C12H24O2	123232-57-5	10



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ClientSampleId :
UI-1

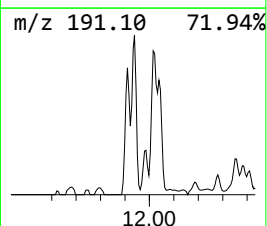
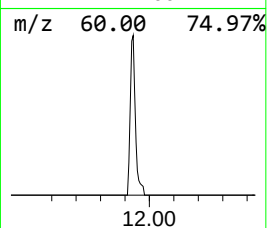
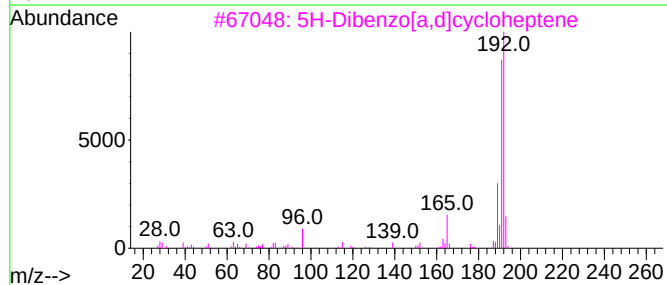
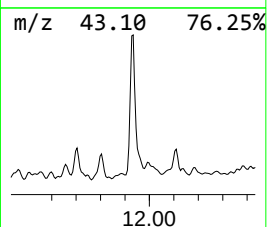
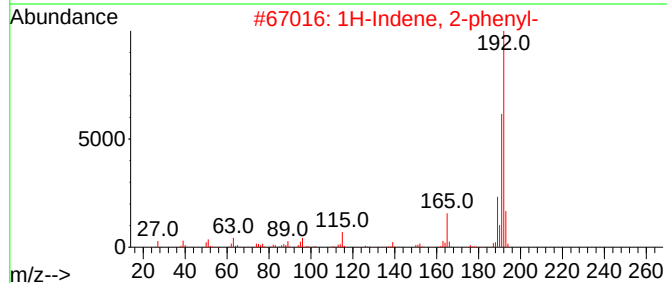
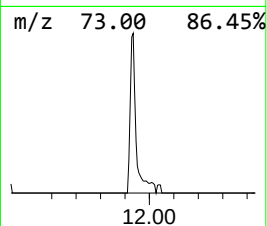
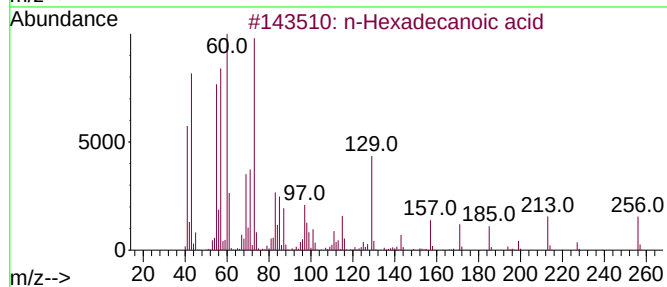
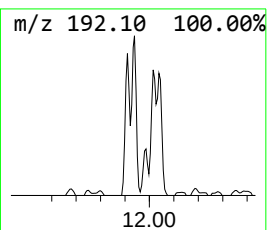
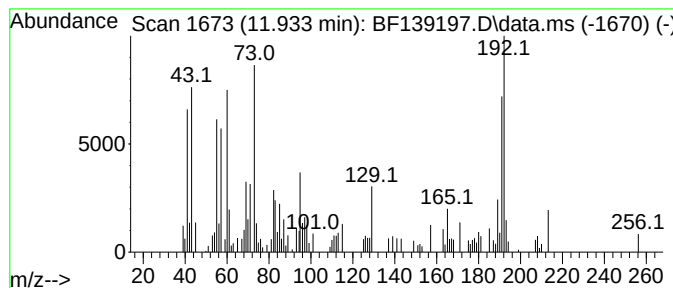
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 n-Hexadecanoic acid Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.933	2.73 ng	96913	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	96
2			1H-Indene, 2-phenyl-	192	C15H12	004505-48-0	83
3			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	80
4			Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	47
5			Naphtho[1,2-b]norbornadiene	192	C15H12	1000210-14-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
Data File : BF139197.D
Acq On : 27 Aug 2024 20:28
Operator : RC/JU
Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

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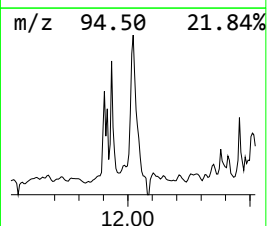
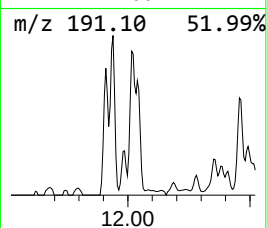
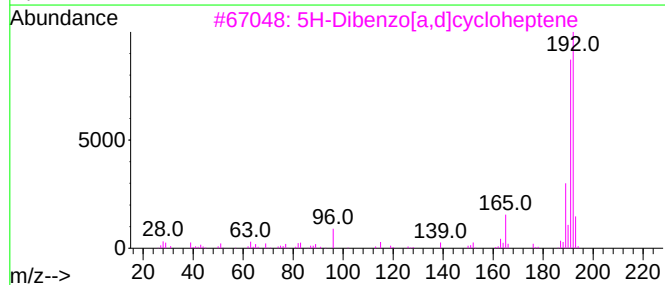
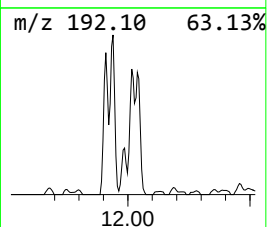
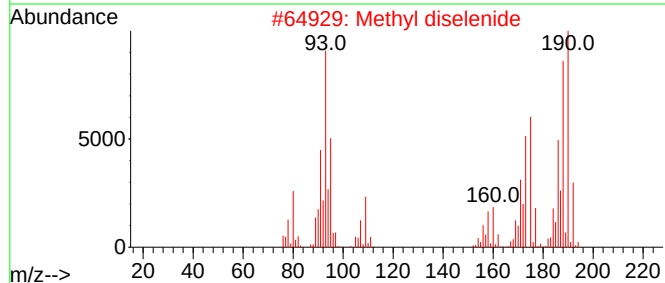
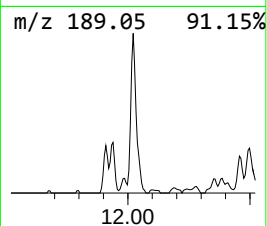
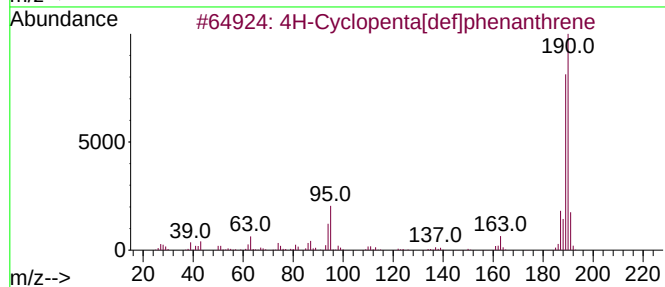
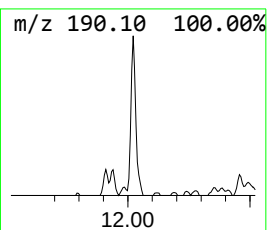
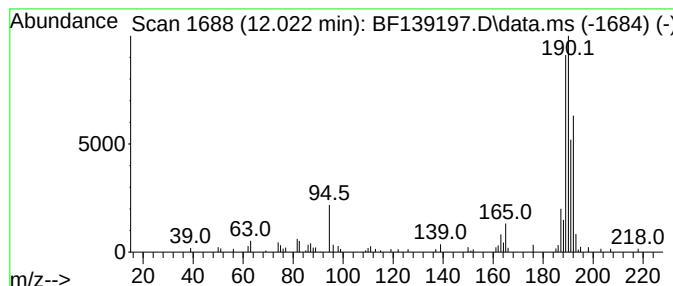
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.022	3.27 ng	115882	Phenanthrene-d10	11.410

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	60
2			Methyl diselenide	190	C2H6Se2	007101-31-7	43
3			5H-Dibenzo[a,d]cycloheptene	192	C15H12	000256-81-5	35
4			Cyclohexanone, 2-(2-furanylmethy...	190	C12H14O2	056053-07-7	27
5			Benzene, 1,1'-(1,2-propadienyld...	192	C15H12	014251-57-1	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
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Acq On : 27 Aug 2024 20:28
Operator : RC/JU
Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

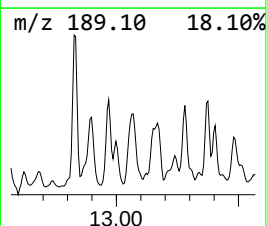
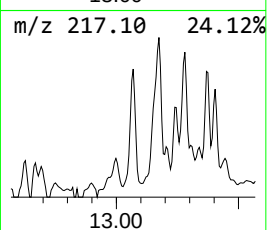
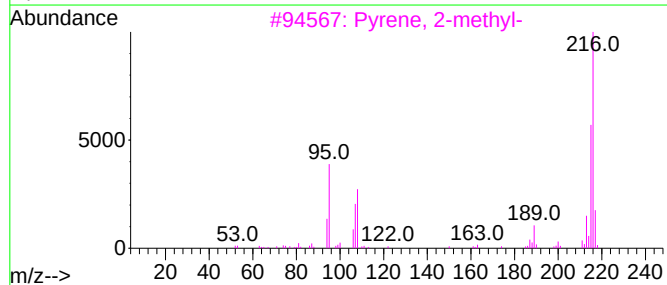
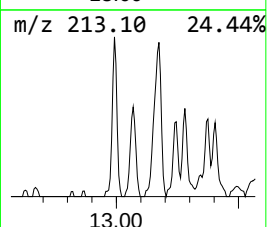
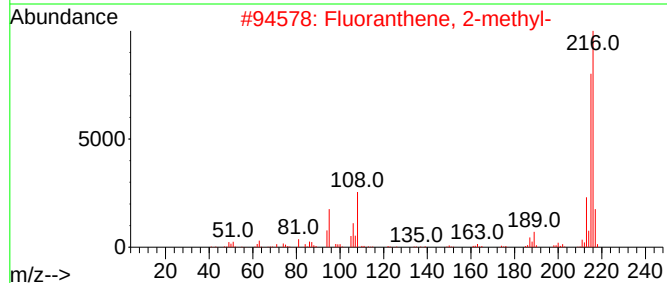
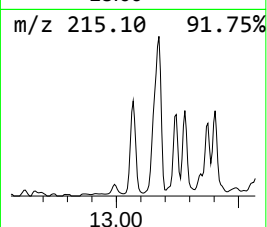
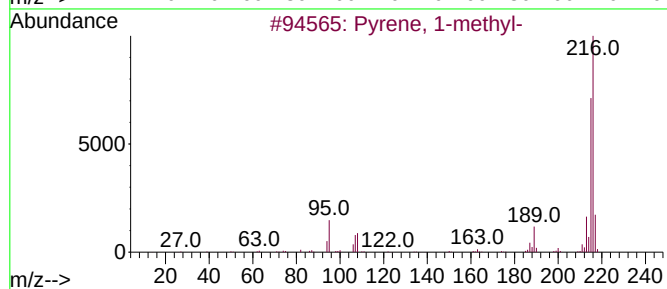
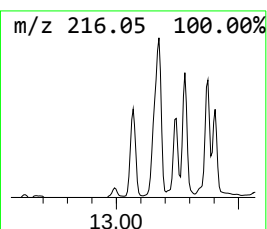
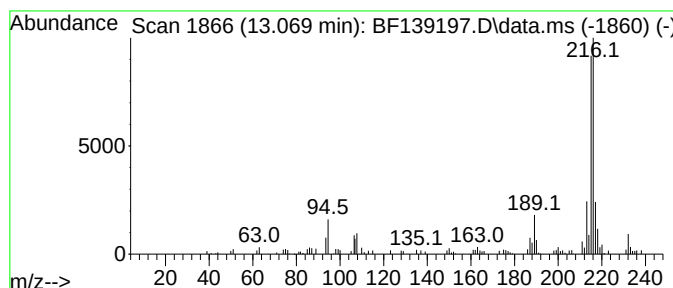
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Pyrene, 1-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.069	2.21 ng	109228	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	86
4	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81
5	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
Data File : BF139197.D
Acq On : 27 Aug 2024 20:28
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Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

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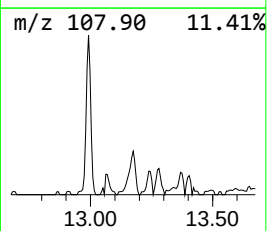
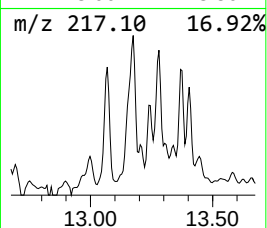
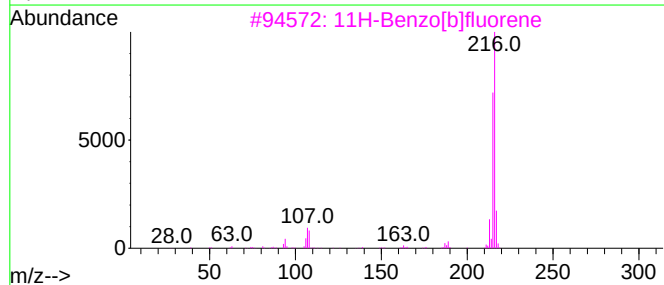
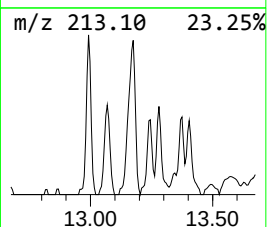
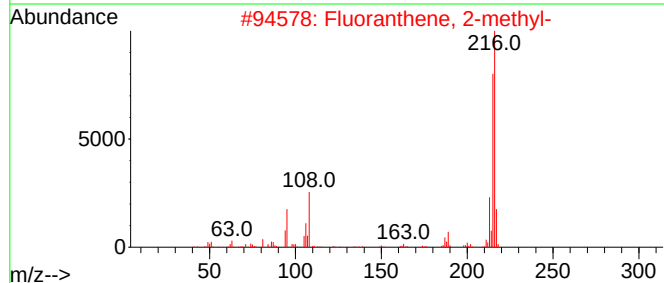
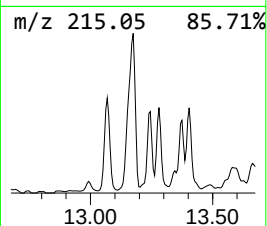
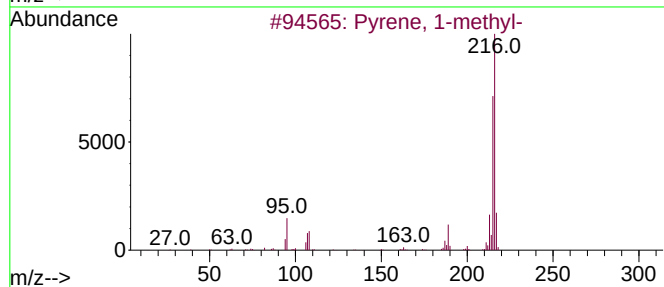
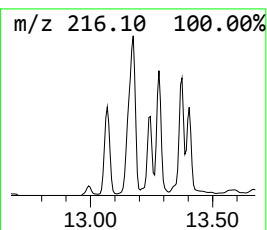
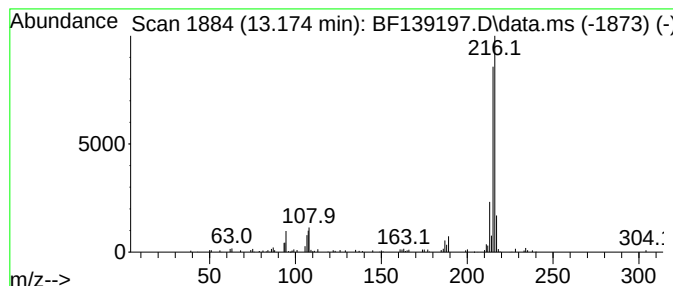
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Fluoranthene, 2-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.174	2.90 ng	143759	Chrysene-d12	14.045

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
2			Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	90
3			11H-Benzo[b]fluorene	216	C17H12	000243-17-4	90
4			11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5			7H-Benzo[c]fluorene	216	C17H12	000205-12-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
Data File : BF139197.D
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Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

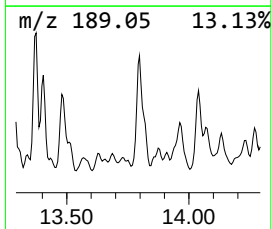
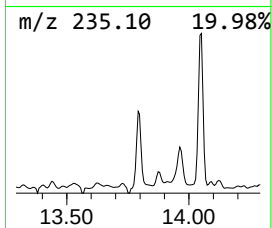
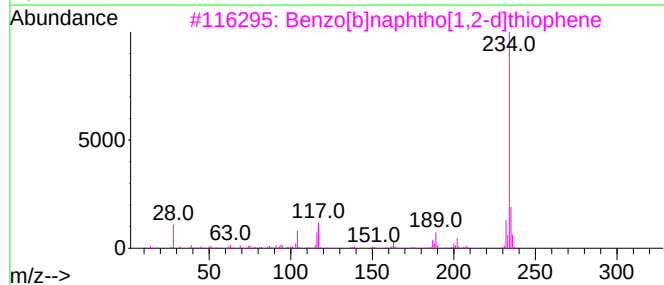
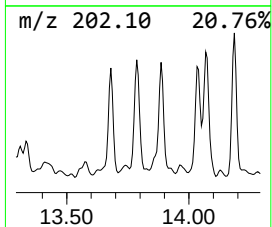
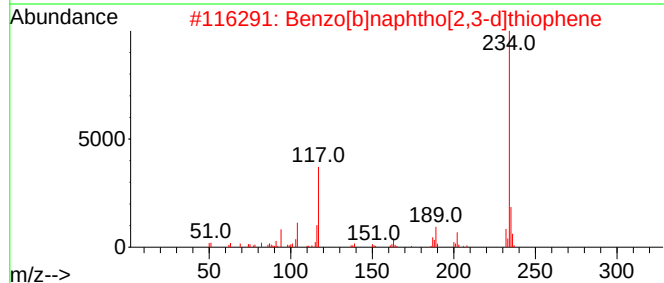
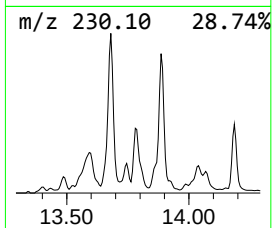
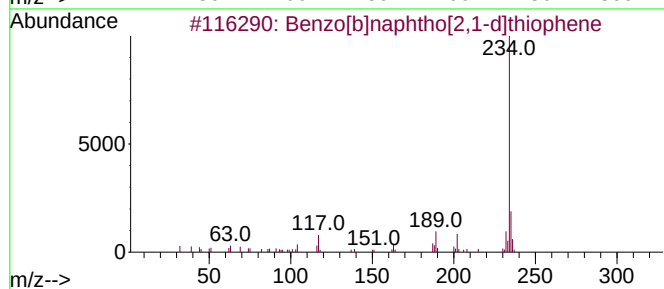
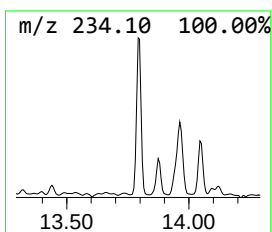
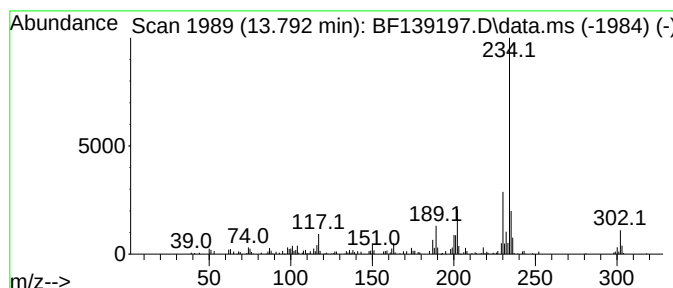
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.792	2.00 ng	99159	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	70
3	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	64
4	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	58
5	pyrido[1',2':1,2]imidazo[4,5-b]q...	234	C14H10N4	1000401-06-3	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
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Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

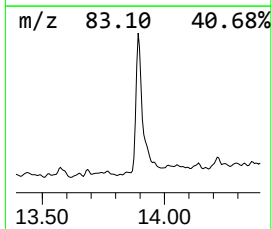
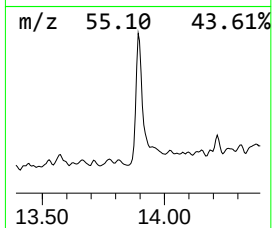
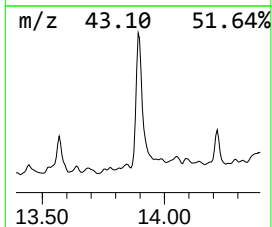
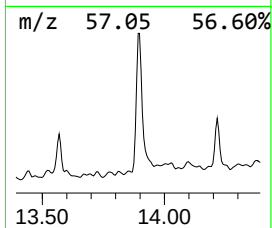
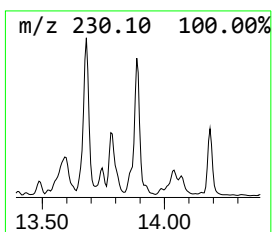
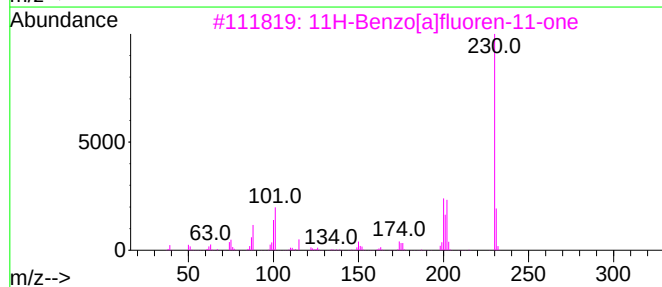
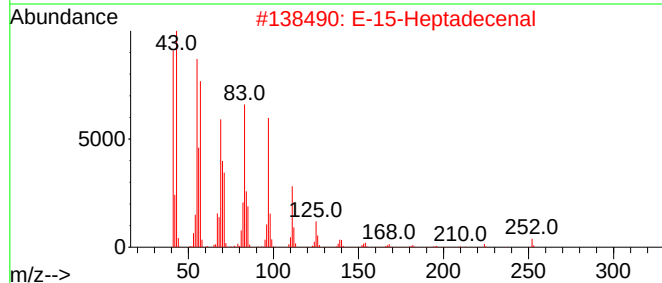
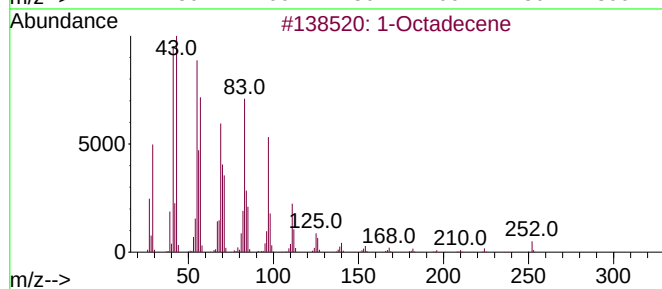
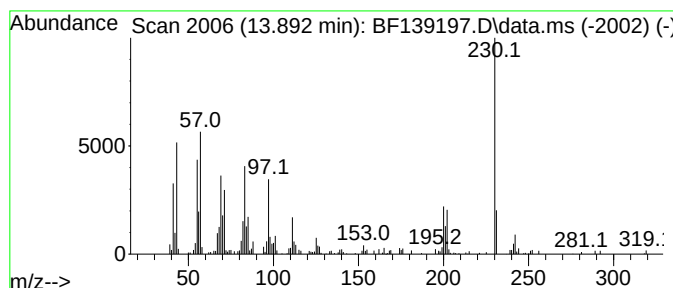
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Octadecene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.892	2.78 ng	137607	Chrysene-d12	14.045	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Octadecene	252	C18H36	000112-88-9	91
2	E-15-Heptadecenal	252	C17H32O	1000130-97-9	90
3	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	60
4	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	55
5	Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	48



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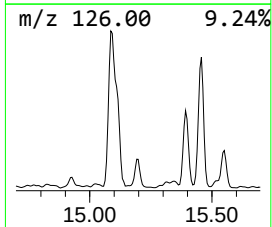
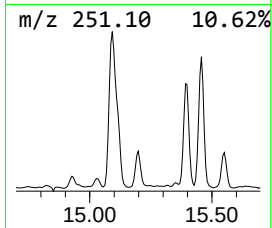
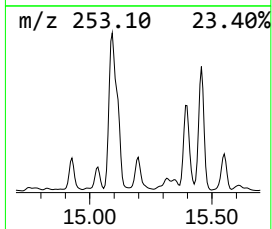
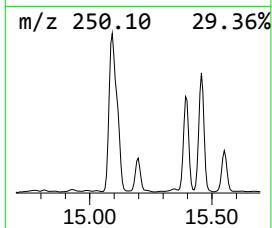
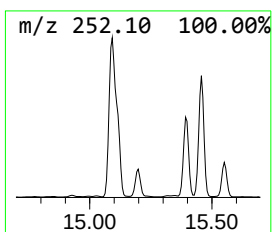
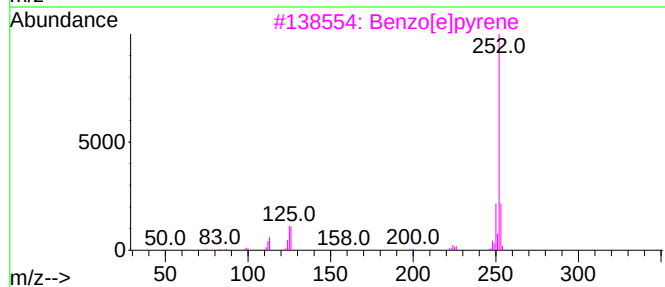
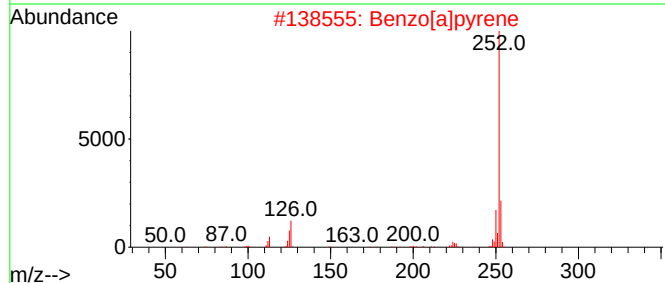
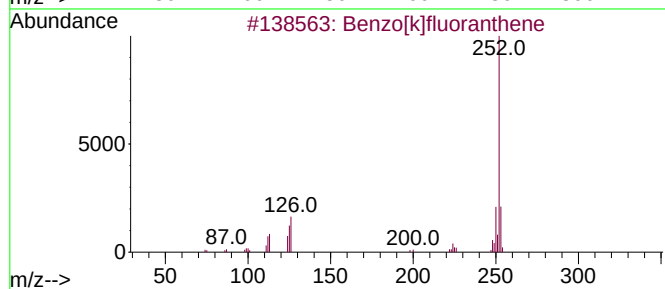
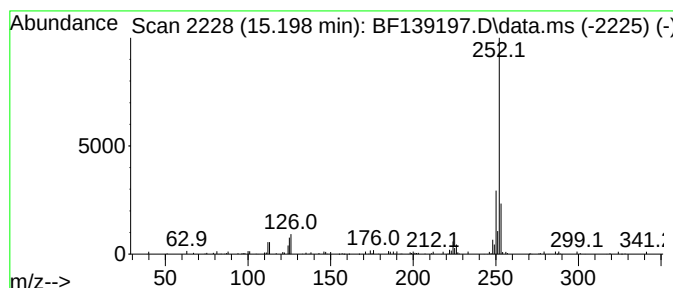
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.198	3.60 ng	65370	Perylene-d12	15.521	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[k]fluoranthene	252	C20H12	000207-08-9	90
2	Benzo[a]pyrene	252	C20H12	000050-32-8	89
3	Benzo[e]pyrene	252	C20H12	000192-97-2	81
4	Benzo[b]fluoranthene	252	C20H12	000205-99-2	81
5	(1-Cyclopentenyl)ferrocene	252	C15H16Fe	012260-67-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
Data File : BF139197.D
Acq On : 27 Aug 2024 20:28
Operator : RC/JU
Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
UI-1

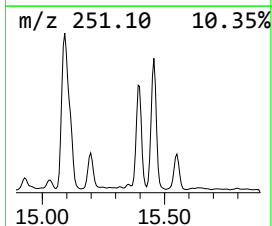
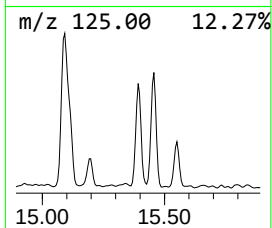
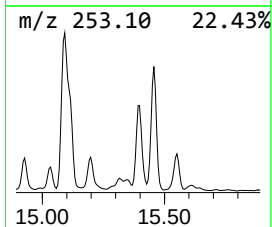
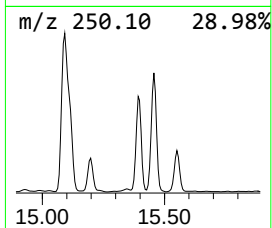
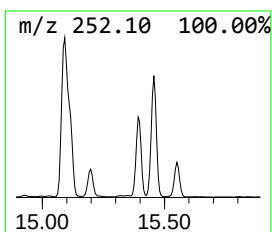
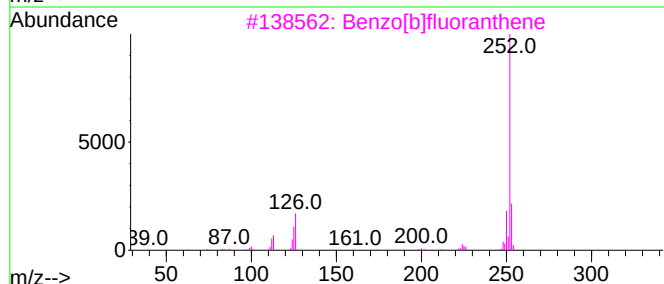
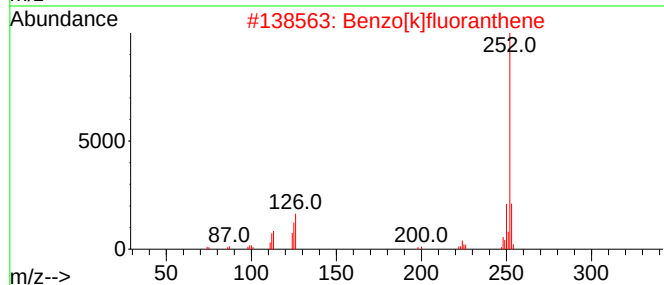
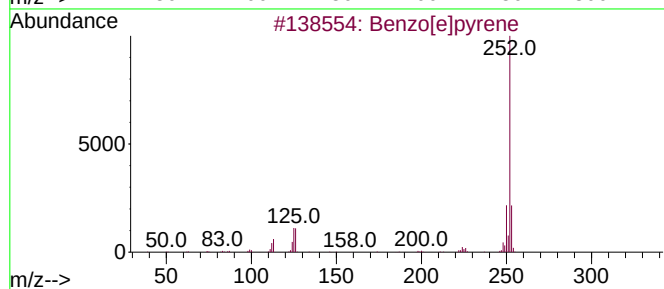
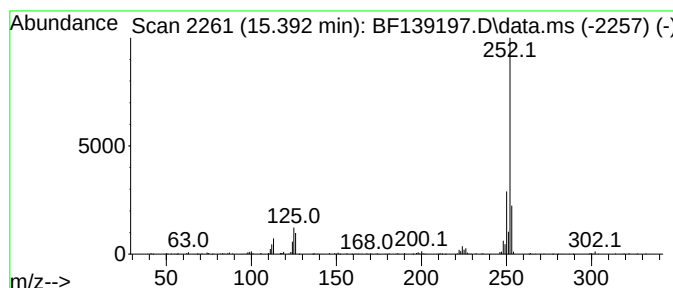
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzo[j]fluoranthene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.392	11.38 ng	206640	Perylene-d12	15.521

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	94
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	93
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	93
5			Benzo[j]fluoranthene	252	C20H12	000205-82-3	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF082724\
Data File : BF139197.D
Acq On : 27 Aug 2024 20:28
Operator : RC/JU
Sample : P3732-01
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
UI-1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF082724.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...	2.181	60.2	ng	1647120	1	6.892	547318	20.0
2-Pentanone, 4-...	5.104	3.9	ng	106265	1	6.892	547318	20.0
1-Phenoxypropan...	8.481	2.3	ng	77214	2	8.175	674206	20.0
unknown10.022	10.022	27.6	ng	931066	3	9.922	675702	20.0
n-Hexadecanoic ...	11.933	2.7	ng	96913	4	11.410	709439	20.0
4H-Cyclopenta[d...	12.022	3.3	ng	115882	4	11.410	709439	20.0
Pyrene, 1-methyl-	13.069	2.2	ng	109228	5	14.045	990258	20.0
Fluoranthene, 2...	13.174	2.9	ng	143759	5	14.045	990258	20.0
Benzo[b]naphtho...	13.792	2.0	ng	99159	5	14.045	990258	20.0
1-Octadecene	13.892	2.8	ng	137607	5	14.045	990258	20.0
Benzo[e]pyrene	15.198	3.6	ng	65370	6	15.521	363228	20.0
Benzo[j]fluoran...	15.392	11.4	ng	206640	6	15.521	363228	20.0