

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
 Data File : BF138701.D
 Acq On : 31 Jul 2024 14:30
 Operator : RC/JU
 Sample : P3376-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-6-Full Waste Class

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-BF073024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF138701.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.352	209	214	223	rBV	7470	23581	1.77%	0.211%
2	5.057	499	504	514	rVB	49886	75260	5.66%	0.673%
3	5.469	568	574	592	rVB	817700	1073533	80.75%	9.602%
4	6.481	740	746	762	rBV	802176	1045664	78.65%	9.352%
5	6.675	775	779	792	rBV	11640	20231	1.52%	0.181%
6	6.846	803	808	813	rBV	249946	310385	23.35%	2.776%
7	7.404	897	903	908	rBV	545358	676271	50.87%	6.049%
8	7.657	942	946	952	rBV	17541	23040	1.73%	0.206%
9	8.122	1020	1025	1034	rVB	294834	368766	27.74%	3.298%
10	8.440	1074	1079	1093	rBV	41594	68170	5.13%	0.610%
11	9.198	1202	1208	1213	rBV	911942	1117339	84.04%	9.994%
12	9.875	1318	1323	1328	rBV	294372	366575	27.57%	3.279%
13	9.969	1334	1339	1345	rBV	25444	38364	2.89%	0.343%
14	10.663	1452	1457	1473	rVB	469515	631639	47.51%	5.649%
15	10.792	1473	1479	1484	rBV3	11846	18767	1.41%	0.168%
16	10.845	1484	1488	1491	rBV	17528	23337	1.76%	0.209%
17	10.881	1491	1494	1497	rBV2	15518	17568	1.32%	0.157%
18	10.992	1507	1513	1515	rVV2	12791	28023	2.11%	0.251%
19	11.028	1515	1519	1522	rVV3	27717	41954	3.16%	0.375%
20	11.063	1522	1525	1529	rVV	24065	37289	2.80%	0.334%
21	11.104	1529	1532	1540	rVB	17128	25460	1.91%	0.228%
22	11.222	1548	1552	1560	rVB7	9756	13812	1.04%	0.124%
23	11.363	1571	1576	1584	rBV2	314458	436693	32.85%	3.906%
24	11.892	1662	1666	1672	rVB	72179	100794	7.58%	0.902%
25	11.975	1677	1680	1690	rVB2	15931	29926	2.25%	0.268%
26	12.575	1777	1782	1787	rBV	139862	188862	14.21%	1.689%
27	12.675	1795	1799	1804	rBV3	17738	27308	2.05%	0.244%
28	12.804	1812	1821	1825	rBV	127634	199778	15.03%	1.787%
29	12.839	1825	1827	1835	rVB3	21401	34795	2.62%	0.311%
30	12.951	1840	1846	1852	rVB	1034148	1329520	100.00%	11.891%
31	13.027	1854	1859	1866	rBV7	10905	22333	1.68%	0.200%
32	13.133	1870	1877	1880	rVV	17370	31597	2.38%	0.283%
33	13.239	1891	1895	1902	rVB3	14727	22800	1.71%	0.204%
34	13.427	1923	1927	1936	rVB	50831	104456	7.86%	0.934%
35	13.504	1937	1940	1946	rBV2	19437	28290	2.13%	0.253%
36	13.639	1960	1963	1967	rBV	14317	19637	1.48%	0.176%

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Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	13.751	1979	1982	1985	rBV2	13514	19632	1.48%	0.176%
38	13.857	1995	2000	2009	rBV3	55763	116791	8.78%	1.045%
39	14.004	2016	2025	2035	rBV2	386221	747792	56.25%	6.688%
40	14.110	2039	2043	2050	rBV3	31250	73368	5.52%	0.656%
41	14.363	2081	2086	2089	rBV	75976	108097	8.13%	0.967%
42	14.469	2101	2104	2106	rBV2	17204	22441	1.69%	0.201%
43	14.757	2148	2153	2160	rBV	201120	297073	22.34%	2.657%
44	14.916	2177	2180	2185	rVB	51377	66617	5.01%	0.596%
45	15.045	2197	2202	2209	rBV	79319	177457	13.35%	1.587%
46	15.104	2209	2212	2216	rVV2	32925	46296	3.48%	0.414%
47	15.163	2217	2222	2237	rVB5	42650	128291	9.65%	1.147%
48	15.345	2249	2253	2259	rVB	42137	62670	4.71%	0.561%
49	15.404	2259	2263	2268	rVV	53625	85195	6.41%	0.762%
50	15.469	2269	2274	2288	rVB	225814	394335	29.66%	3.527%
51	15.686	2308	2311	2318	rVB2	24386	35084	2.64%	0.314%
52	15.916	2345	2350	2357	rBV	47011	82178	6.18%	0.735%
53	16.927	2518	2522	2529	rBV2	22238	46633	3.51%	0.417%
54	17.374	2593	2598	2606	rVB	21895	48870	3.68%	0.437%

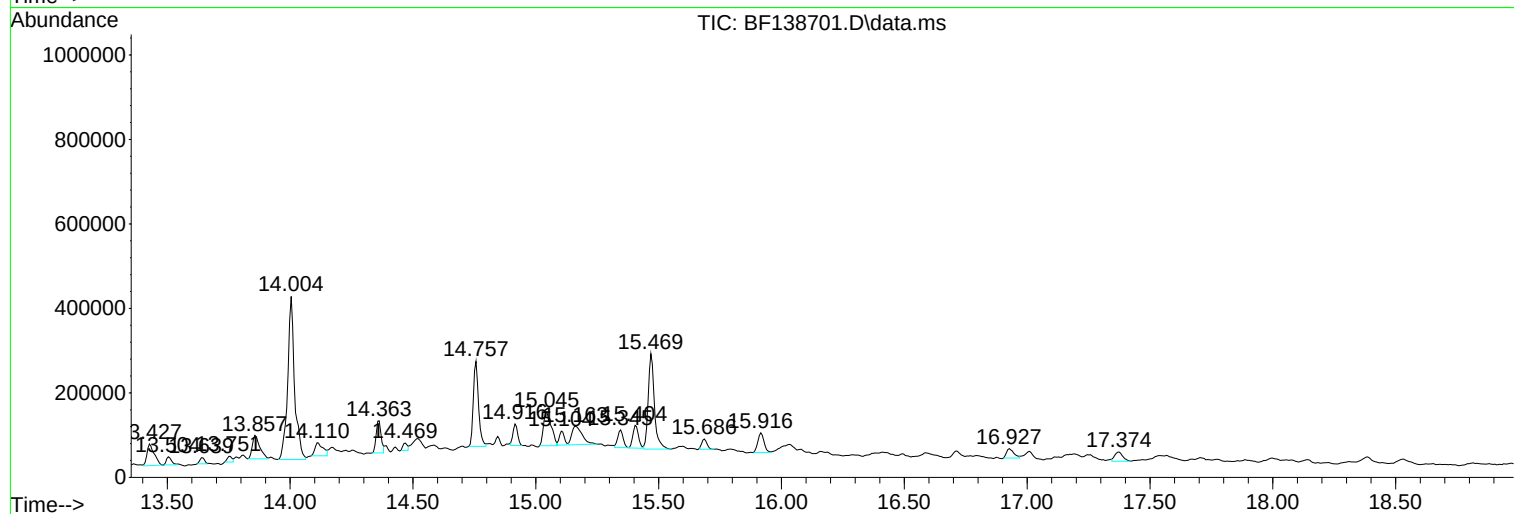
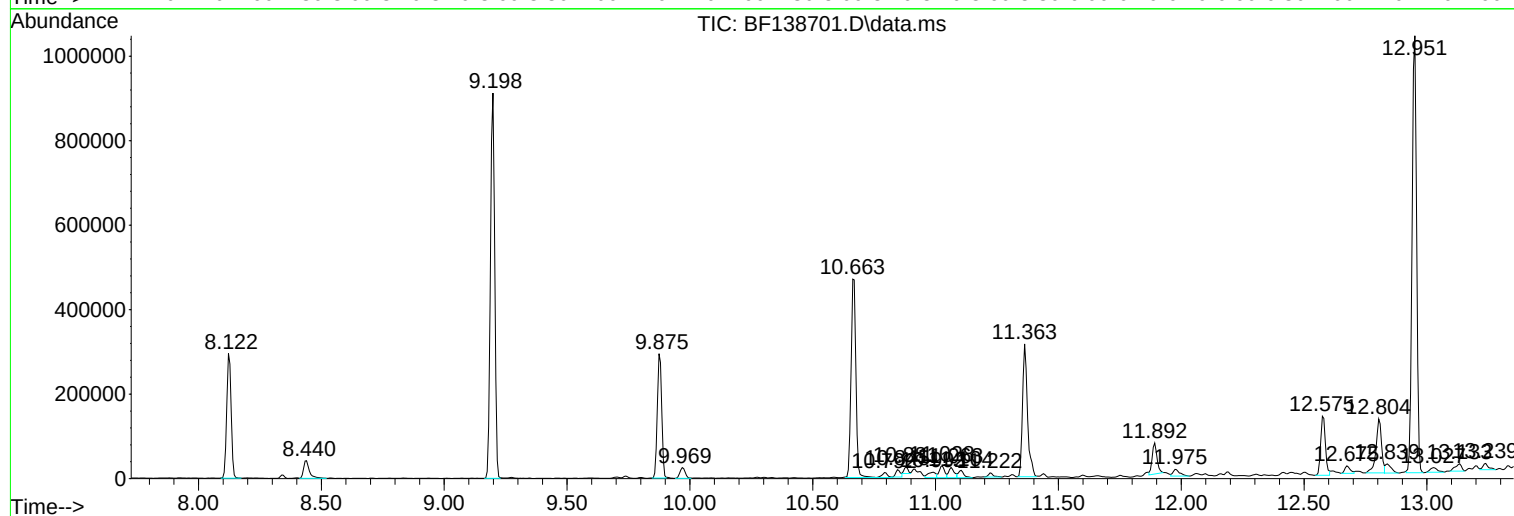
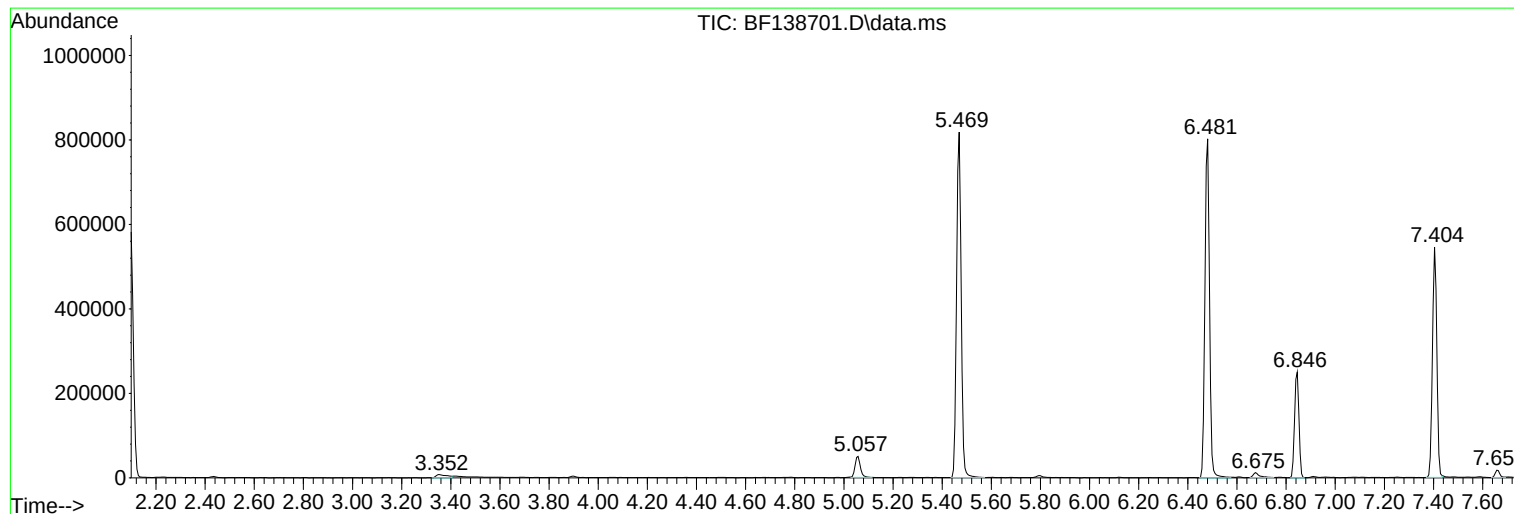
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
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ClientSampleId :
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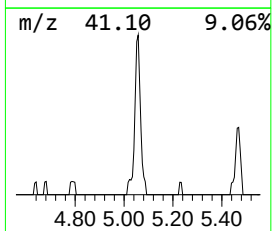
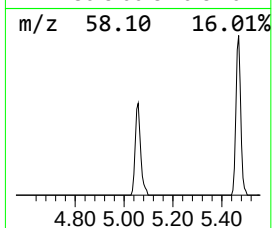
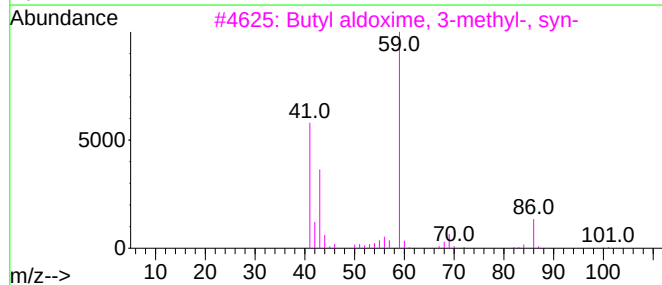
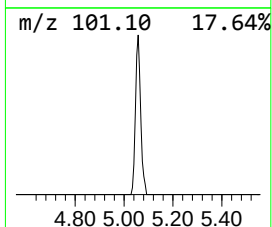
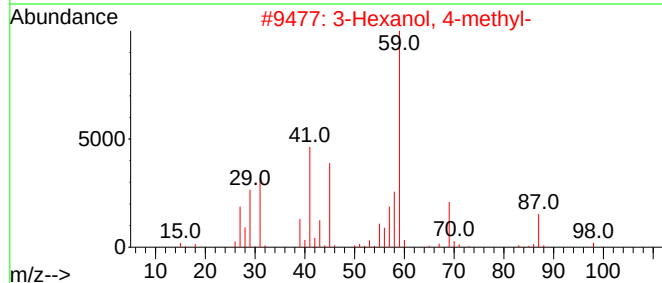
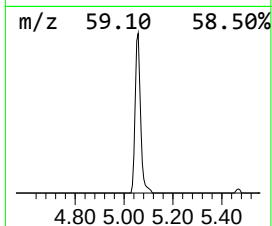
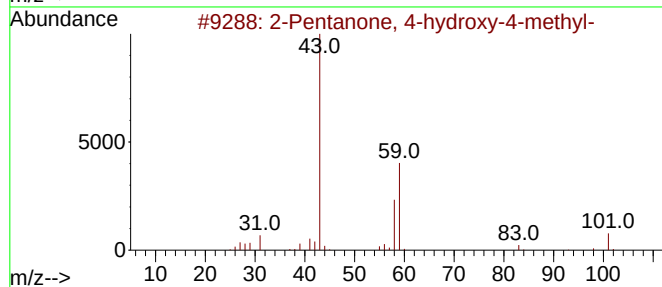
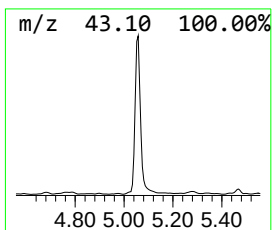
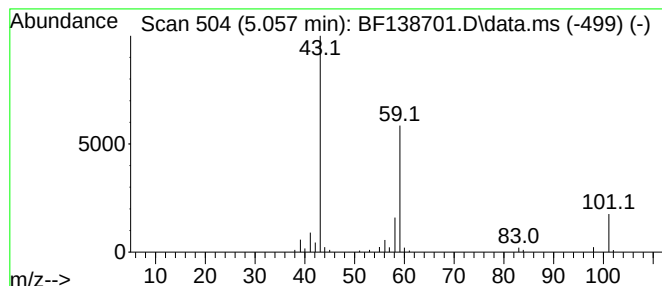
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.057	4.85 ng	75260	1,4-Dichlorobenzene-d4	6.846

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	56		
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33		
3	Butyl aldoxime, 3-methyl-, syn-	101	C5H11NO	005780-40-5	32		
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16		
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12		



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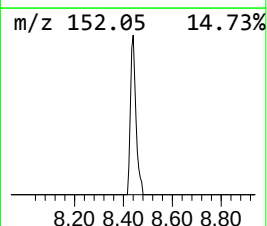
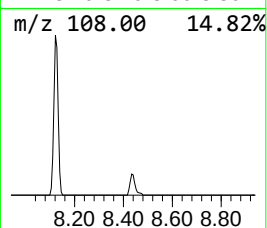
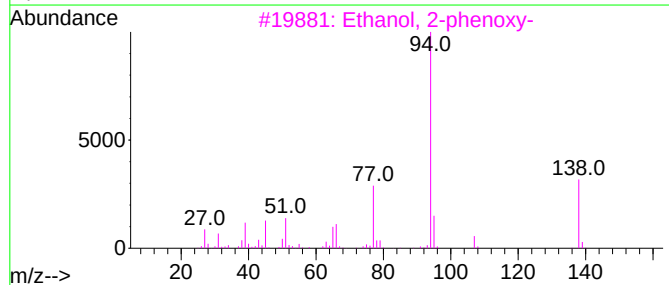
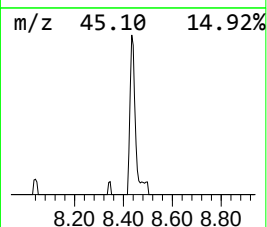
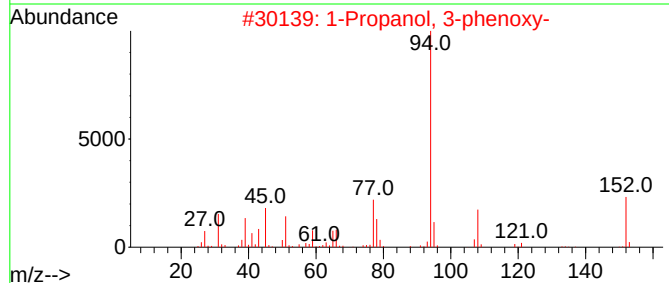
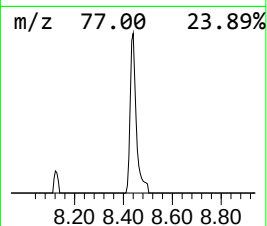
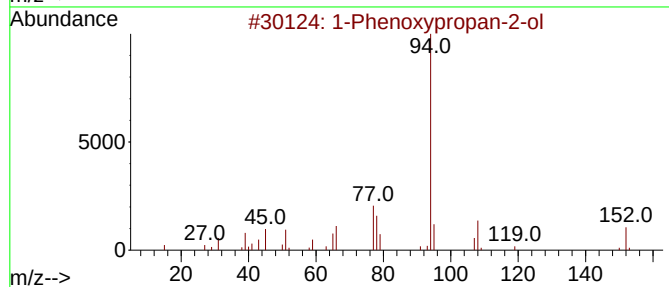
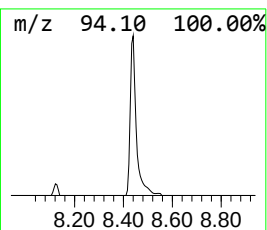
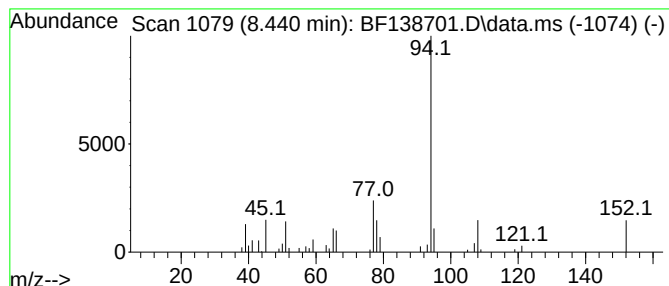
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.440	3.70 ng	68170	Naphthalene-d8	8.122

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	93
2		1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	91
3		Ethanol, 2-phenoxy-	138	C8H10O2	000122-99-6	64
4		1-Propanol, 2-phenoxy-	152	C9H12O2	004169-04-4	59
5		2-Hexen-4-yne, 2-methyl-	94	C7H10	058275-93-7	53



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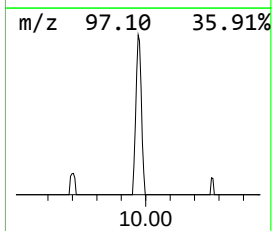
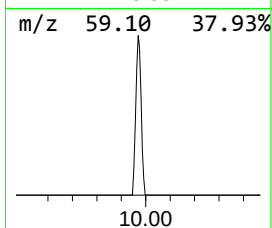
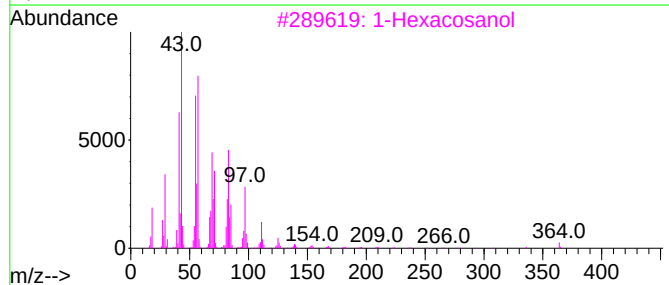
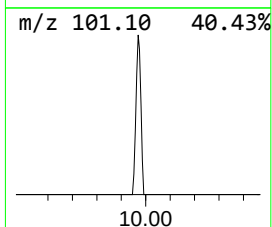
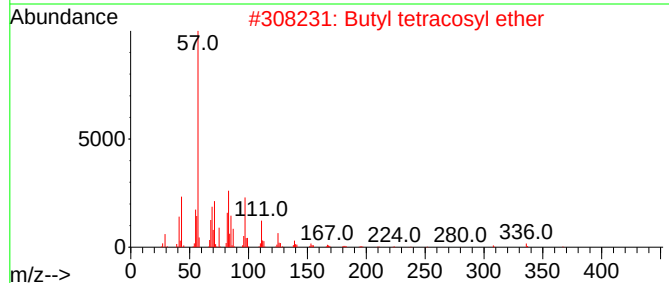
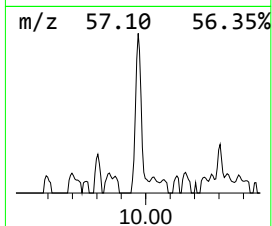
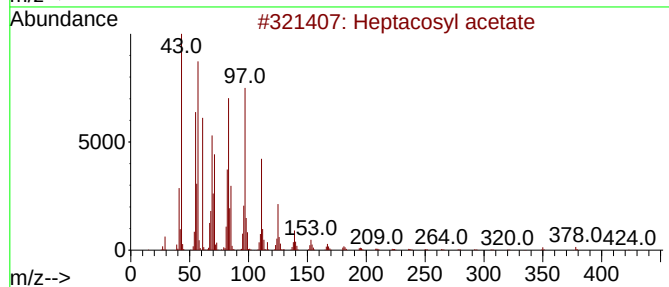
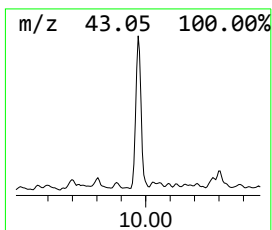
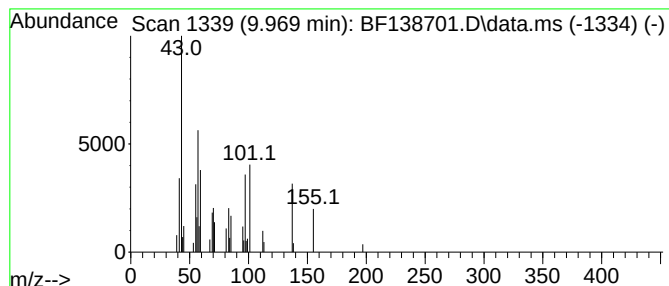
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 unknown9.969 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.969	2.09 ng	38364	Acenaphthene-d10	9.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosyl acetate	438	C29H58O2	1000351-78-2	14
2			Butyl tetracosyl ether	410	C28H58O	1000406-40-9	10
3			1-Hexacosanol	382	C26H54O	000506-52-5	10
4			1-Tetracosene	336	C24H48	010192-32-2	10
5			Carbonic acid, prop-1-en-2-yl te...	298	C18H34O3	1000382-90-9	10



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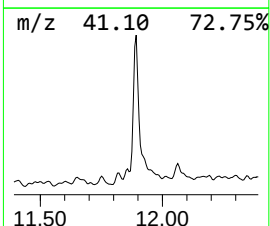
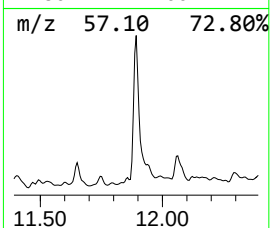
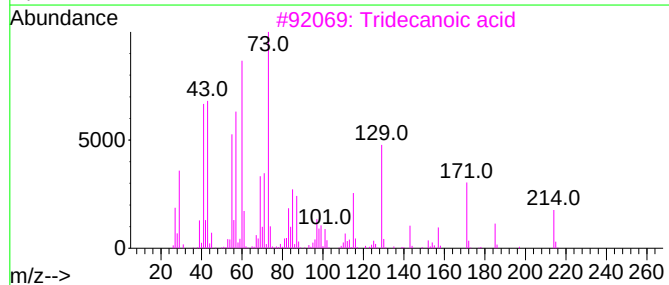
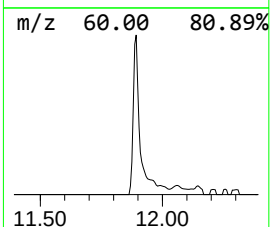
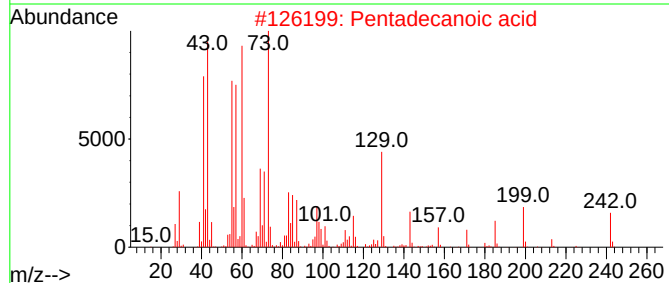
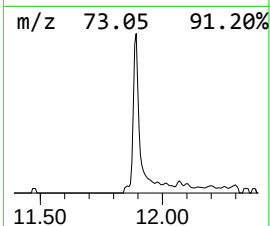
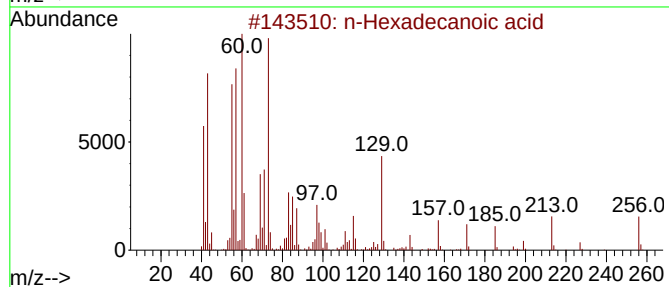
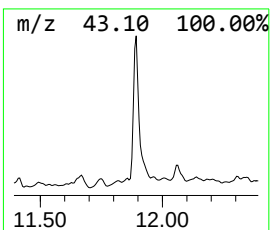
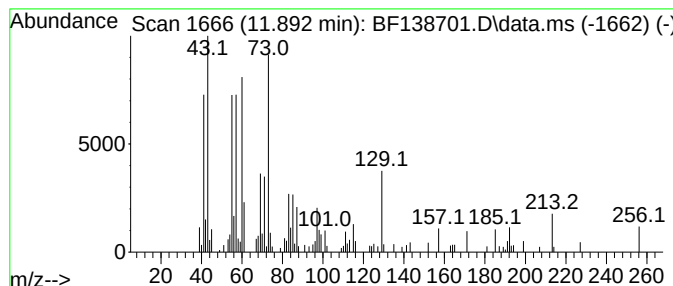
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.892	4.62 ng	100794	Phenanthrene-d10	11.363

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			Pentadecanoic acid	242	C15H30O2	001002-84-2	87
3			Tridecanoic acid	214	C13H26O2	000638-53-9	81
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			Undecanoic acid	186	C11H22O2	000112-37-8	53



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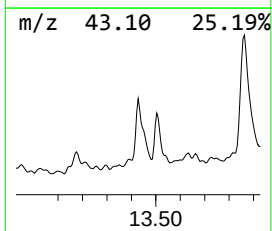
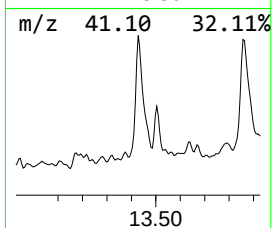
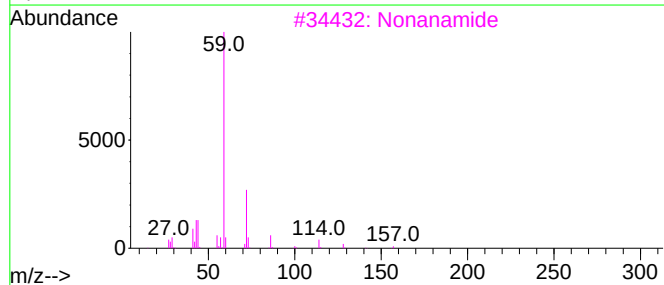
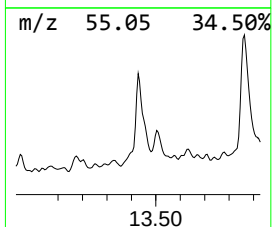
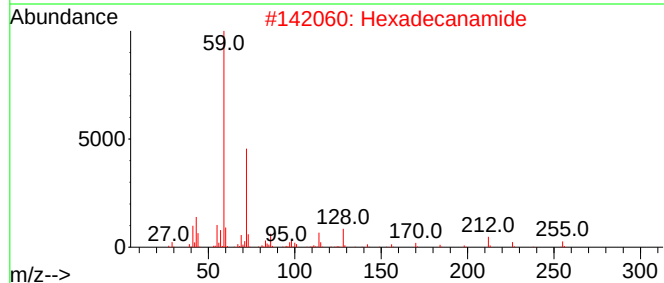
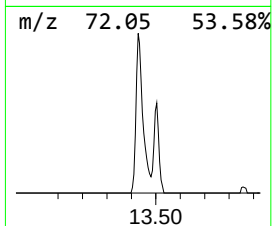
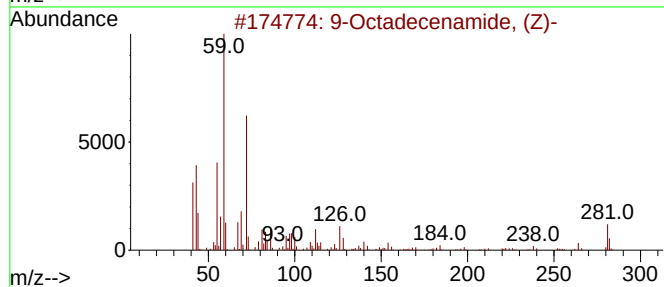
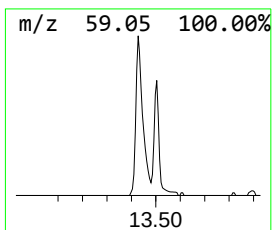
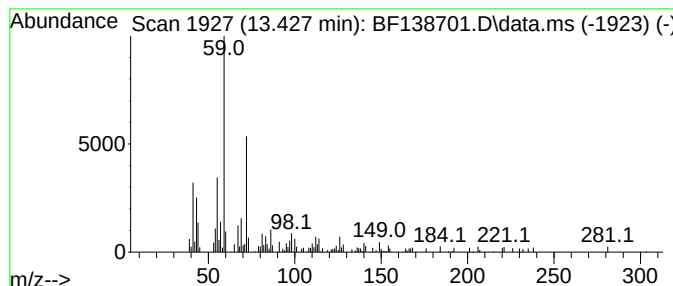
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ClientSampleId :
WC-6-Full Waste Class

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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 9-Octadecenamide, (Z)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.428	2.79 ng	104456	Chrysene-d12	14.004	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	90
2	Hexadecanamide	255	C16H33NO	000629-54-9	78
3	Nonanamide	157	C9H19NO	001120-07-6	64
4	Dodecanamide	199	C12H25NO	001120-16-7	64
5	Octanamide	143	C8H17NO	000629-01-6	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

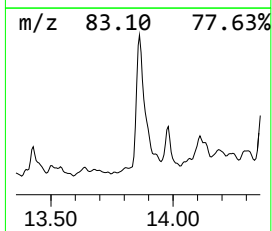
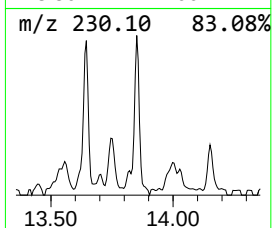
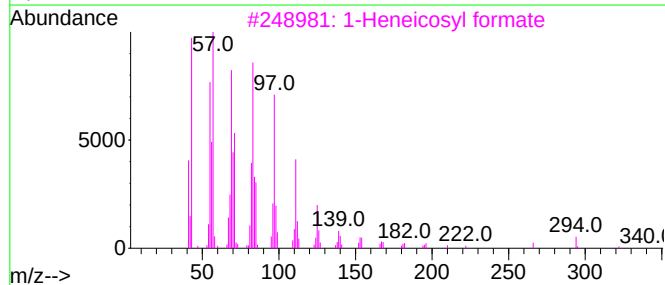
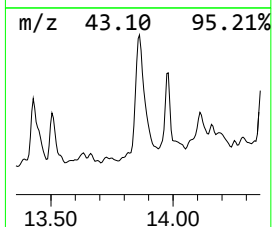
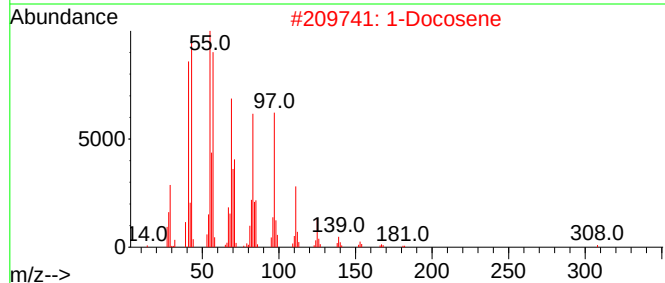
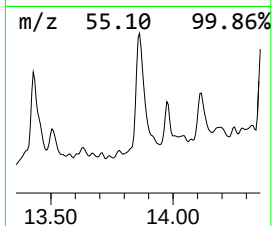
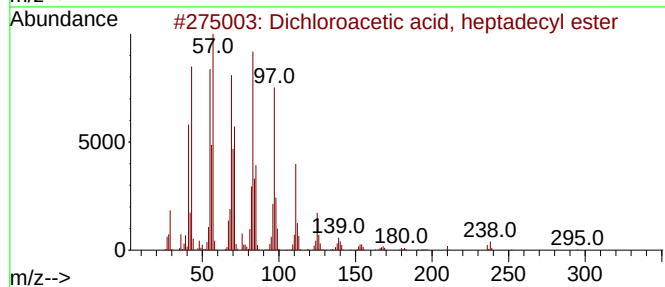
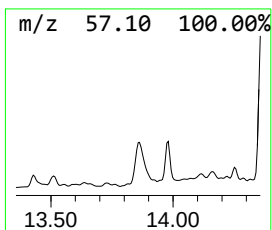
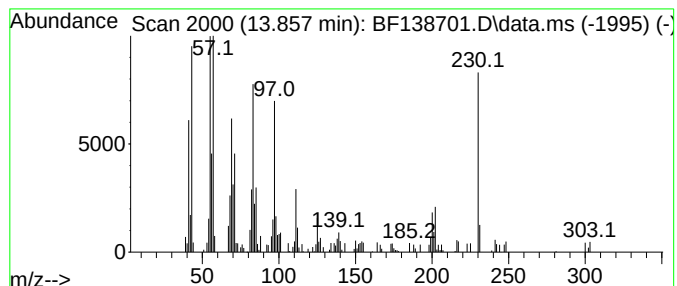
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Dichloroacetic acid, heptad... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.857	3.12 ng	116791	Chrysene-d12	14.004

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	93
2		1-Docosene	308	C22H44	001599-67-3	86
3		1-Heneicosyl formate	340	C22H44O2	077899-03-7	70
4		Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	70
5		1-Hexacosene	364	C26H52	018835-33-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

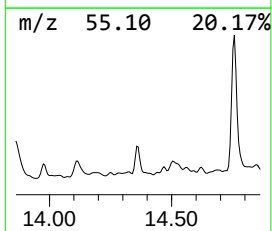
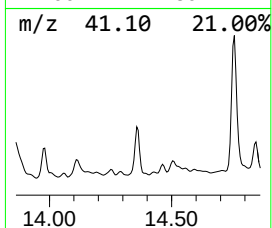
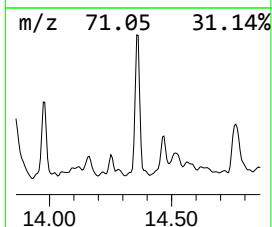
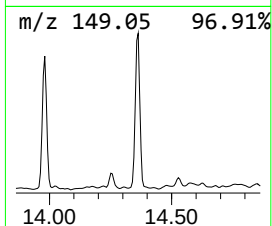
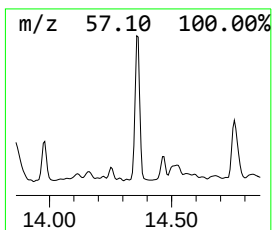
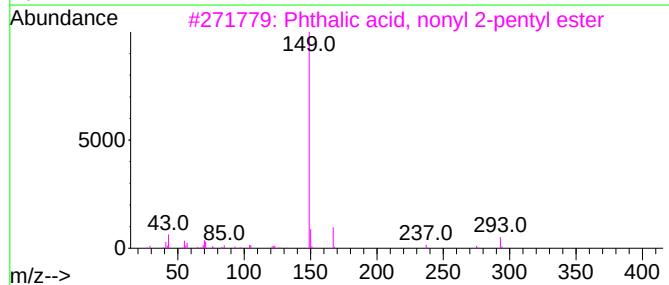
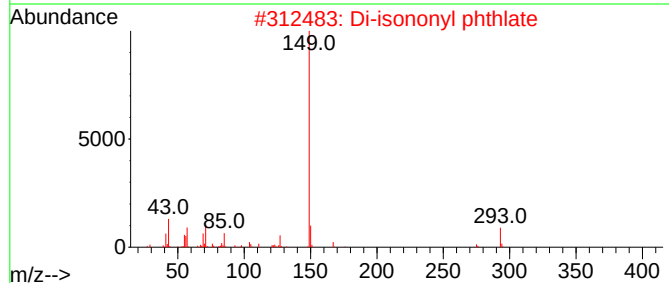
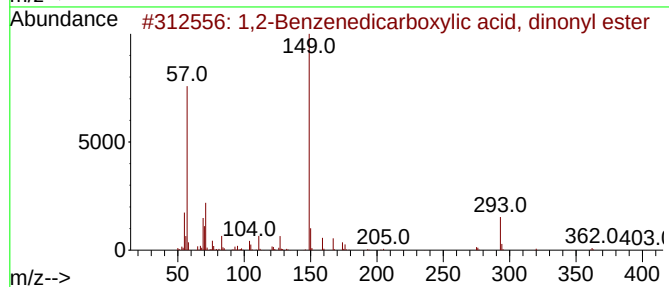
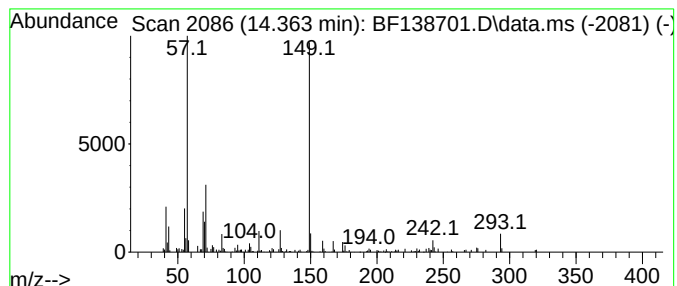
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 1,2-Benzenedicarboxylic aci... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.363	2.89 ng	108097	Chrysene-d12	14.004

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,2-Benzenedicarboxylic acid, di...	418	C26H42O4	000084-76-4	72
2			Di-isononyl phthlate	418	C26H42O4	020548-62-3	53
3			Phthalic acid, nonyl 2-pentyl ester	362	C22H34O4	1000315-48-1	47
4			Phthalic acid, 3-methylbut-3-eny...	360	C22H32O4	1010357-10-7	47
5			1,2-Benzenedicarboxylic acid, di...	362	C22H34O4	003648-21-3	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

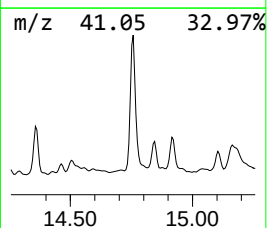
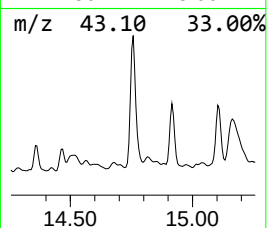
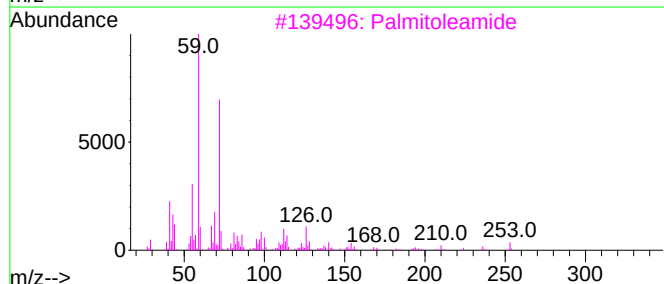
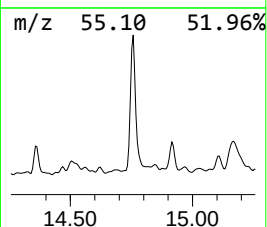
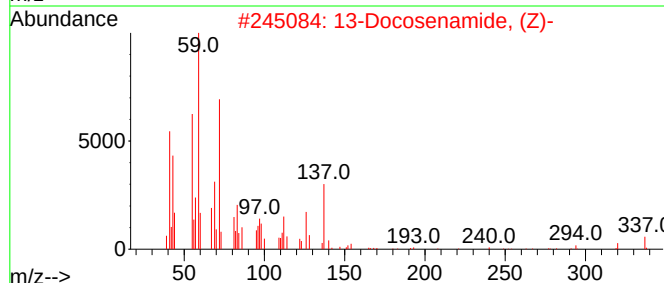
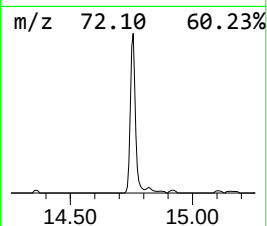
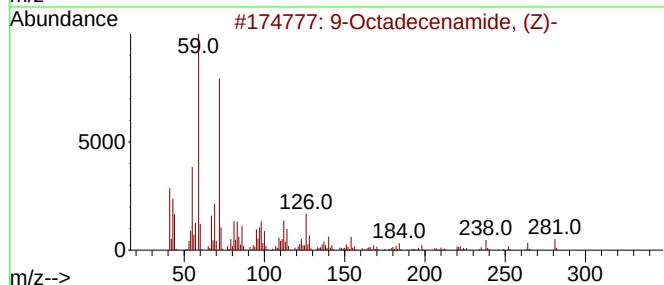
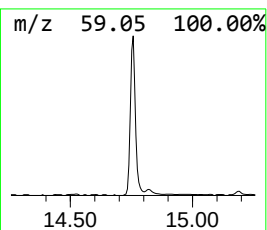
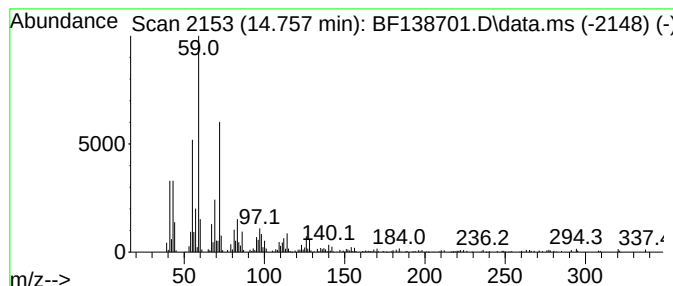
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 13-Docosenamide, (Z)- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.757	15.07 ng	297073	Perylene-d12	15.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenamide, (Z)-			281	C18H35NO	000301-02-0	93
2	13-Docosenamide, (Z)-			337	C22H43NO	000112-84-5	93
3	Palmitoleamide			253	C16H31NO	106010-22-4	72
4	Octadecanamide			283	C18H37NO	000124-26-5	59
5	Butanamide, 3-methyl-			101	C5H11NO	000541-46-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
WC-6-Full Waste Class

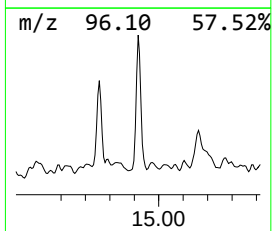
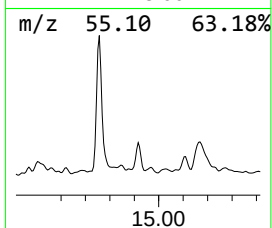
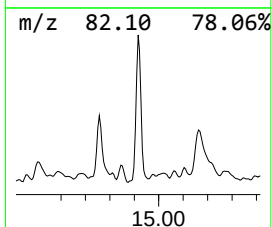
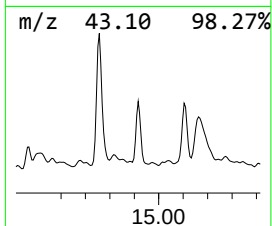
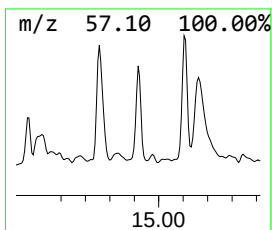
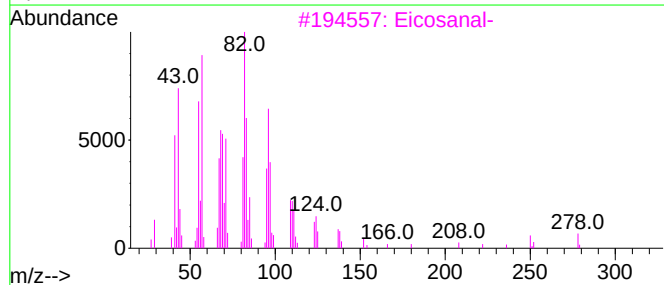
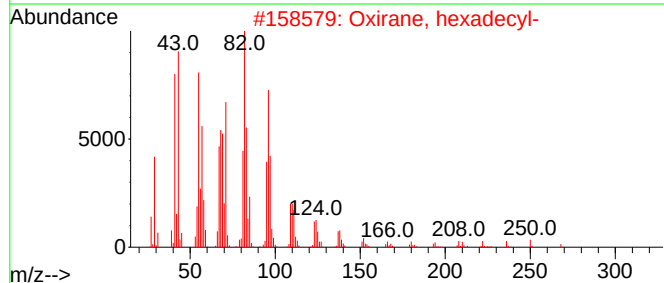
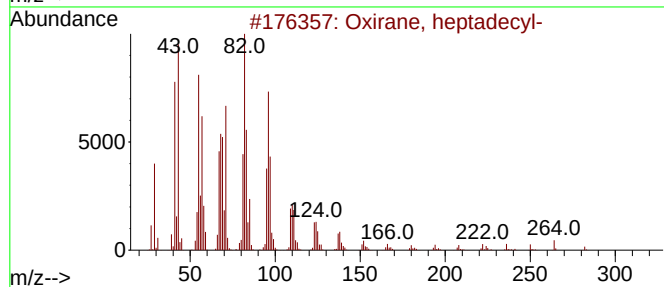
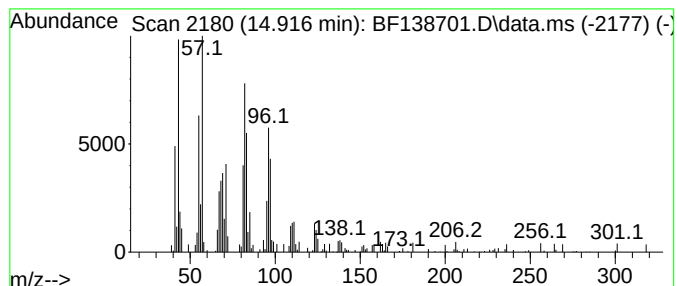
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Oxirane, heptadecyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.916	3.38 ng	66617	Perylene-d12	15.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	81
2			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	76
3			Eicosanal-	296	C20H40O	002400-66-0	70
4			1,16-Hexadecanediol	258	C16H34O2	007735-42-4	58
5			1,19-Eicosadiene	278	C20H38	014811-95-1	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

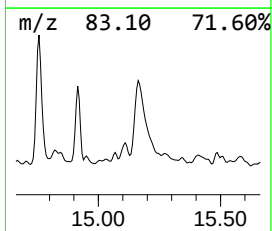
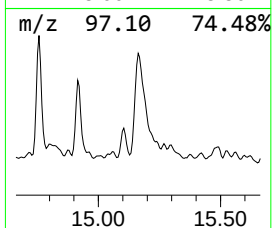
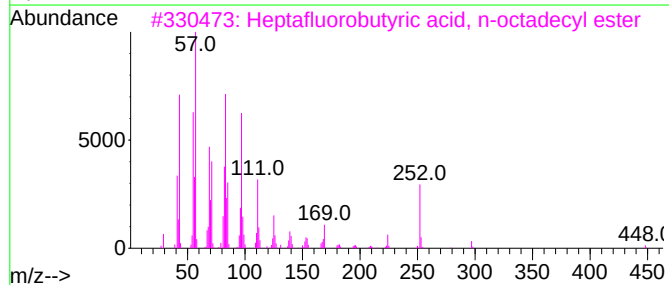
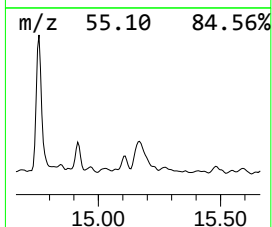
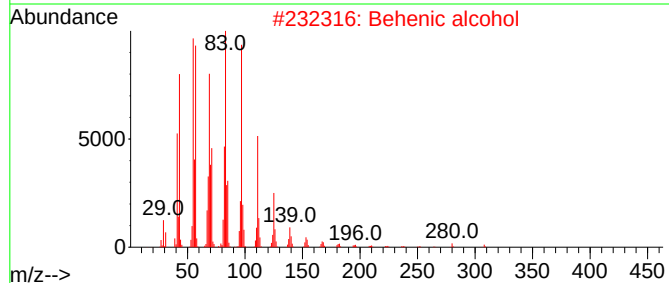
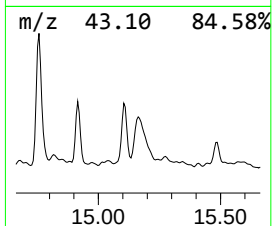
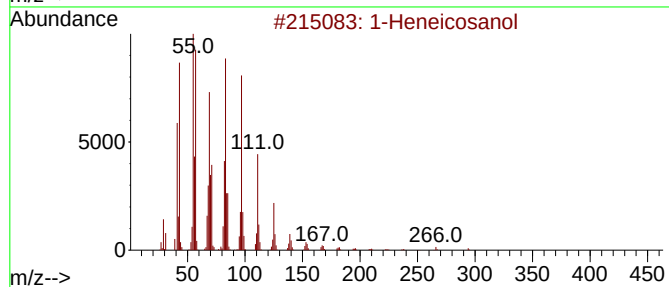
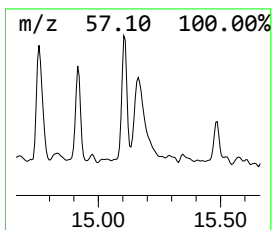
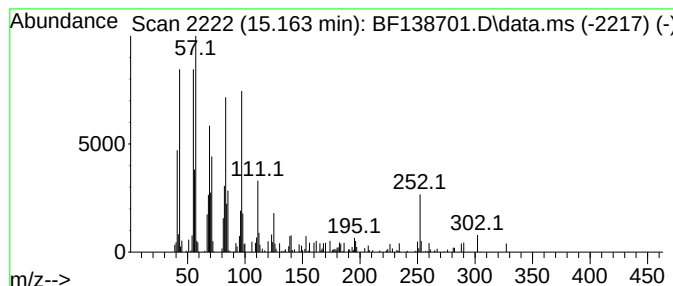
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 1-Heneicosanol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.163	6.51 ng	128291	Perylene-d12	15.469

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	93
2		Behenic alcohol	326	C22H46O	000661-19-8	93
3		Heptafluorobutyric acid, n-octadecyl ester	466	C22H37F7O2	000400-57-7	91
4		Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	91
5		Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

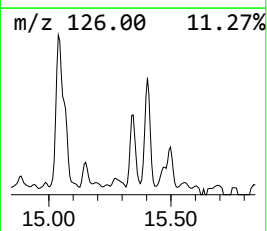
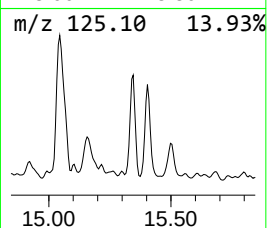
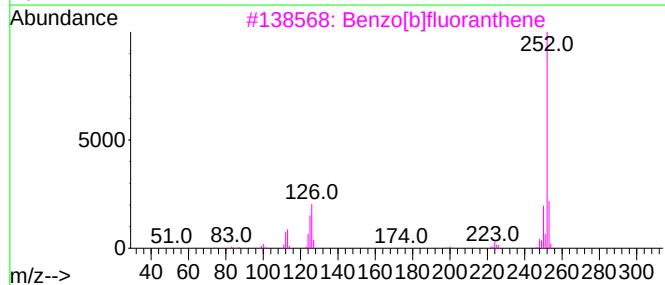
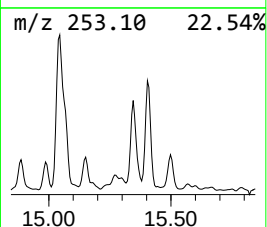
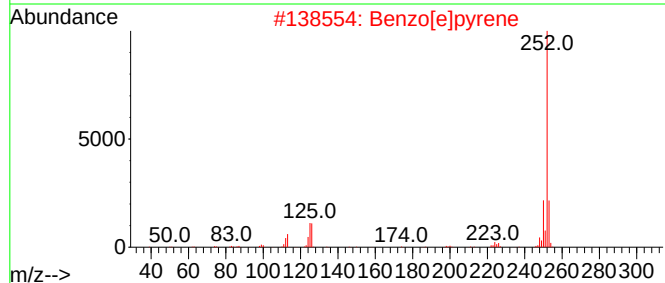
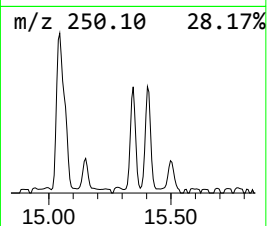
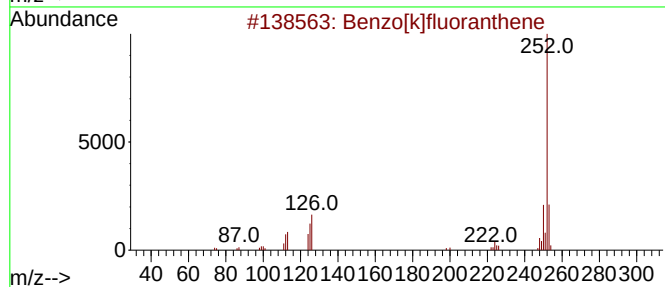
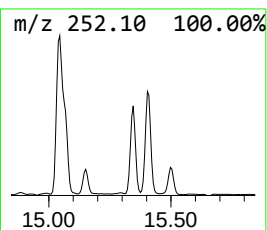
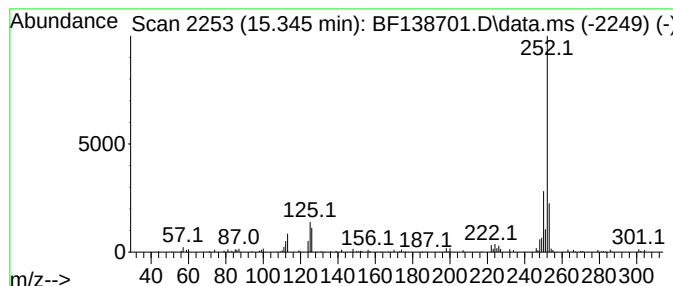
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Benzo[e]pyrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.345	3.18 ng	62670	Perylene-d12	15.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
2			Benzo[e]pyrene	252	C20H12	000192-97-2	96
3			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
4			Benzo[a]pyrene	252	C20H12	000050-32-8	91
5			Perylene	252	C20H12	000198-55-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

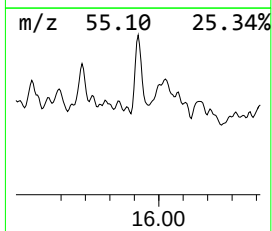
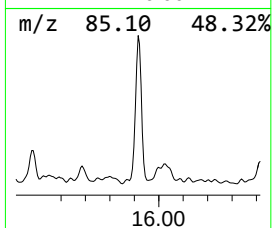
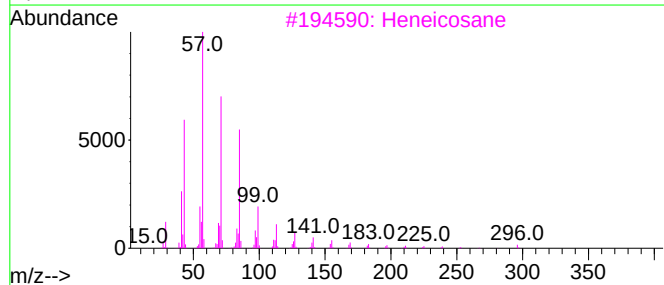
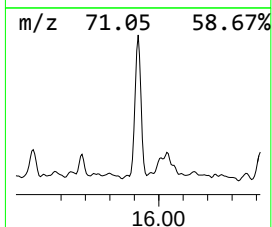
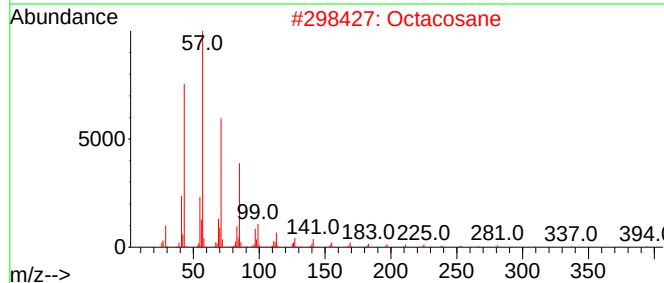
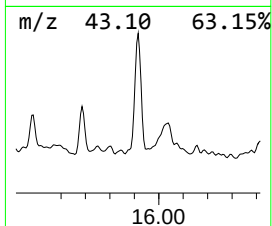
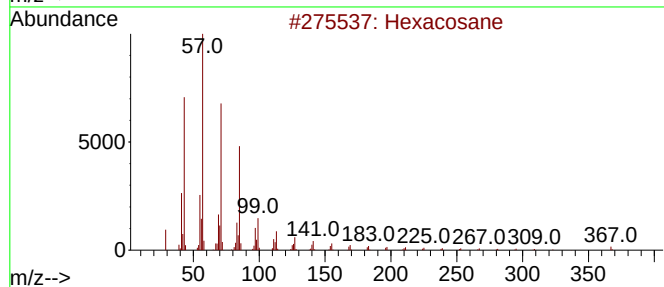
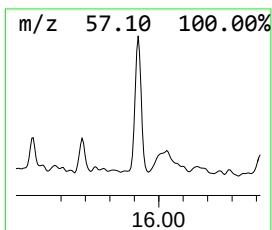
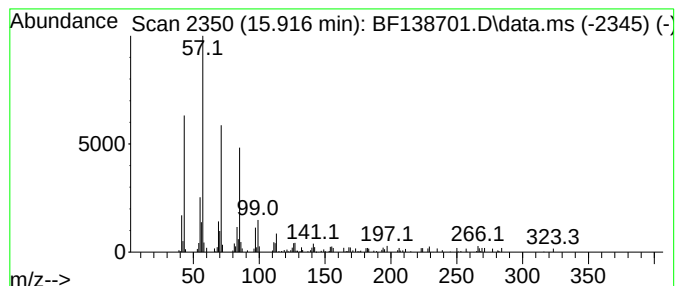
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Hexacosane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.916	4.17 ng	82178	Perylene-d12	15.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexacosane	366	C26H54	000630-01-3	90
2			Octacosane	394	C28H58	000630-02-4	90
3			Heneicosane	296	C21H44	000629-94-7	87
4			Heptadecane	240	C17H36	000629-78-7	87
5			Eicosane	282	C20H42	000112-95-8	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

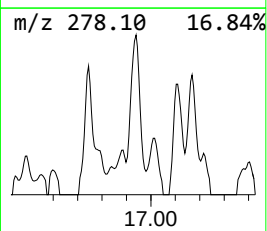
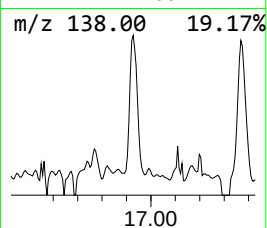
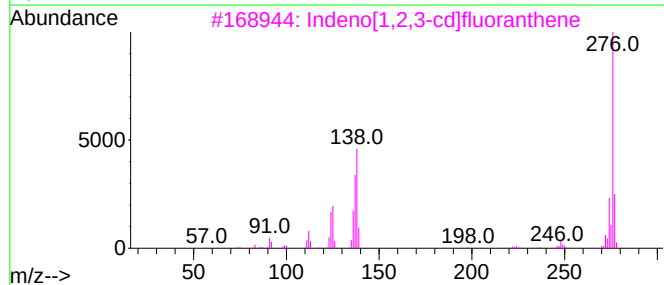
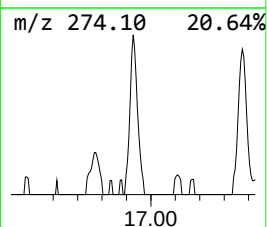
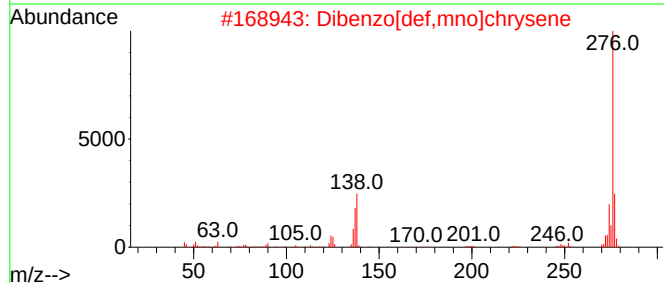
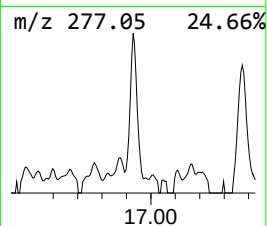
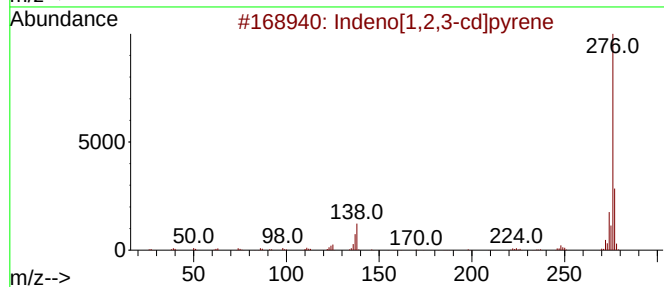
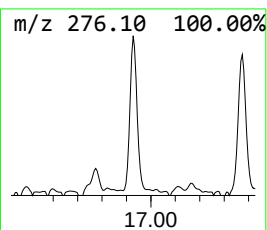
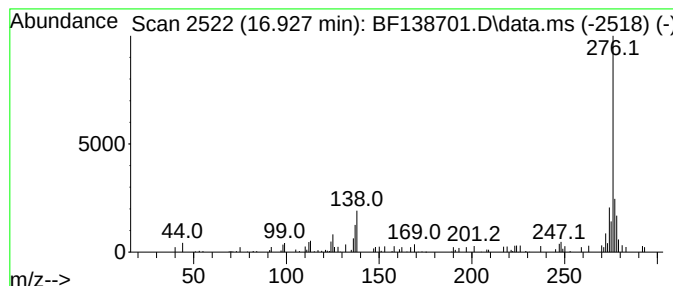
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Dibenzo[def,mno]chrysene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.927	2.37 ng	46633	Perylene-d12	15.469

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indeno[1,2,3-cd]pyrene	276	C22H12	000193-39-5	90
2			Dibenzo[def,mno]chrysene	276	C22H12	000191-26-4	81
3			Indeno[1,2,3-cd]fluoranthene	276	C22H12	000193-43-1	81
4			Benzo[ghi]perylene	276	C22H12	000191-24-2	74
5			2-Thiophenecarbaldehyde, 5-[5-(t...	276	C13H8OS3	1000217-28-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF073124\
Data File : BF138701.D
Acq On : 31 Jul 2024 14:30
Operator : RC/JU
Sample : P3376-04
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-6-Full Waste Class

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF073024.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
2-Pentanone, 4-...	5.057	4.8	ng	75260	1	6.846	310385	20.0
1-Phenoxypropan...	8.440	3.7	ng	68170	2	8.122	368766	20.0
unknown9.969	9.969	2.1	ng	38364	3	9.875	366575	20.0
n-Hexadecanoic ...	11.892	4.6	ng	100794	4	11.363	436693	20.0
9-Octadecenamid...	13.428	2.8	ng	104456	5	14.004	747792	20.0
Dichloroacetic ...	13.857	3.1	ng	116791	5	14.004	747792	20.0
1,2-Benzenedica...	14.363	2.9	ng	108097	5	14.004	747792	20.0
13-Docosenamide...	14.757	15.1	ng	297073	6	15.469	394335	20.0
Oxirane, heptad...	14.916	3.4	ng	66617	6	15.469	394335	20.0
1-Heneicosanol	15.163	6.5	ng	128291	6	15.469	394335	20.0
Benzo[e]pyrene	15.345	3.2	ng	62670	6	15.469	394335	20.0
Hexacosane	15.916	4.2	ng	82178	6	15.469	394335	20.0
Dibenzo[def,mno...	16.927	2.4	ng	46633	6	15.469	394335	20.0