

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
 Data File : BF136922.D
 Acq On : 03 Jan 2024 22:57
 Operator : CG\JU
 Sample : 05899-01
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GHL-SS-08

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-BF122023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF136922.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.169	10	14	19	rVB	15726	19513	1.53%	0.152%
2	3.275	194	202	209	rBV	13719	27689	2.18%	0.215%
3	5.022	495	499	506	rVB	25804	33309	2.62%	0.259%
4	5.381	557	560	565	rBV	16393	17514	1.38%	0.136%
5	5.428	565	568	583	rBV	1230964	1194250	93.90%	9.274%
6	5.634	600	603	606	rBV	16787	15818	1.24%	0.123%
7	6.034	668	671	677	rBV5	11174	16070	1.26%	0.125%
8	6.087	677	680	683	rBV	22893	17534	1.38%	0.136%
9	6.310	715	718	724	rBV2	34303	37278	2.93%	0.289%
10	6.434	736	739	749	rBV	1273808	1157926	91.04%	8.992%
11	6.522	750	754	758	rVB4	18319	27342	2.15%	0.212%
12	6.616	766	770	772	rBV	62539	56872	4.47%	0.442%
13	6.739	789	791	793	rBV2	20382	17198	1.35%	0.134%
14	6.781	795	798	803	rVB2	474299	421856	33.17%	3.276%
15	6.834	803	807	810	rBV2	29548	34056	2.68%	0.264%
16	6.869	810	813	815	rVV3	19891	23586	1.85%	0.183%
17	6.922	819	822	826	rVV2	54511	54235	4.26%	0.421%
18	6.957	826	828	830	rVB	30469	21847	1.72%	0.170%
19	7.004	833	836	838	rVV	24782	18186	1.43%	0.141%
20	7.051	838	844	847	rVB3	60246	81579	6.41%	0.633%
21	7.081	847	849	851	rBV2	44792	31086	2.44%	0.241%
22	7.116	851	855	858	rBV2	51767	57021	4.48%	0.443%
23	7.163	858	863	866	rVV2	63117	71923	5.65%	0.558%
24	7.198	866	869	875	rVB5	22637	30388	2.39%	0.236%
25	7.298	882	886	887	rBV2	38171	37755	2.97%	0.293%
26	7.345	890	894	896	rVV	838332	756016	59.44%	5.871%
27	7.369	896	898	902	rVB	176438	144318	11.35%	1.121%
28	7.422	902	907	909	rVB4	34170	52723	4.15%	0.409%
29	7.551	927	929	931	rBV2	27082	23173	1.82%	0.180%
30	7.734	957	960	963	rVB2	13503	16083	1.26%	0.125%
31	7.804	968	972	975	rBV4	16271	18015	1.42%	0.140%
32	7.975	997	1001	1009	rBV3	12788	27075	2.13%	0.210%
33	8.057	1012	1015	1017	rVV	560418	420796	33.08%	3.268%
34	8.075	1017	1018	1023	rVB	90428	72877	5.73%	0.566%
35	8.769	1133	1136	1141	rBV	26517	22099	1.74%	0.172%
36	8.839	1146	1148	1151	rBV2	15885	19600	1.54%	0.152%

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Integrator: RTE

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Sampling : 1

Start Thrs: 0.2

Stop Thrs : 0

Filtering: 5

Min Area: 3 % of largest Peak

Max Peaks: 100

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-BF122023.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	9.133	1194	1198	1201	rBV	1646628	1271871	100.00%	9.876%
38	9.810	1310	1313	1317	rBV	461485	413922	32.54%	3.214%
39	10.022	1345	1349	1354	rVB	32344	28539	2.24%	0.222%
40	10.363	1404	1407	1414	rVB	71696	60338	4.74%	0.469%
41	10.604	1445	1448	1459	rBV	691077	590616	46.44%	4.586%
42	11.304	1563	1567	1569	rBV	407006	339646	26.70%	2.637%
43	11.327	1569	1571	1575	rVB	231067	175011	13.76%	1.359%
44	11.380	1575	1580	1584	rBV	462464	377853	29.71%	2.934%
45	11.545	1602	1608	1612	rBV	127640	123073	9.68%	0.956%
46	11.845	1655	1659	1665	rBV	377046	413178	32.49%	3.208%
47	12.527	1771	1775	1777	rBV	169273	169648	13.34%	1.317%
48	12.551	1777	1779	1789	rVB3	90225	111058	8.73%	0.862%
49	12.627	1789	1792	1796	rBV	154279	159864	12.57%	1.241%
50	12.751	1809	1813	1816	rVB	124854	131821	10.36%	1.024%
51	12.898	1834	1838	1842	rVB	1051657	932395	73.31%	7.240%
52	13.086	1868	1870	1872	rVB	25975	19642	1.54%	0.153%
53	13.151	1879	1881	1883	rBV	20611	15758	1.24%	0.122%
54	13.616	1957	1960	1965	rVB	200383	177620	13.97%	1.379%
55	13.757	1982	1984	1991	rBV6	23761	37782	2.97%	0.293%
56	13.851	1996	2000	2008	rBV5	48674	105720	8.31%	0.821%
57	13.945	2012	2016	2017	rBV	348851	354264	27.85%	2.751%
58	13.963	2017	2019	2022	rVV	414644	375303	29.51%	2.914%
59	13.986	2022	2023	2027	rVB	82911	57134	4.49%	0.444%
60	14.351	2083	2085	2088	rVB3	29879	28645	2.25%	0.222%
61	14.451	2100	2102	2105	rVB3	36880	27627	2.17%	0.215%
62	14.492	2105	2109	2117	rBV	710031	699587	55.00%	5.432%
63	15.010	2193	2197	2200	rBV	70048	100394	7.89%	0.780%
64	15.315	2246	2249	2252	rVB	32088	32717	2.57%	0.254%
65	15.380	2257	2260	2266	rVB	61955	84658	6.66%	0.657%
66	15.445	2266	2271	2275	rBV	290854	367700	28.91%	2.855%

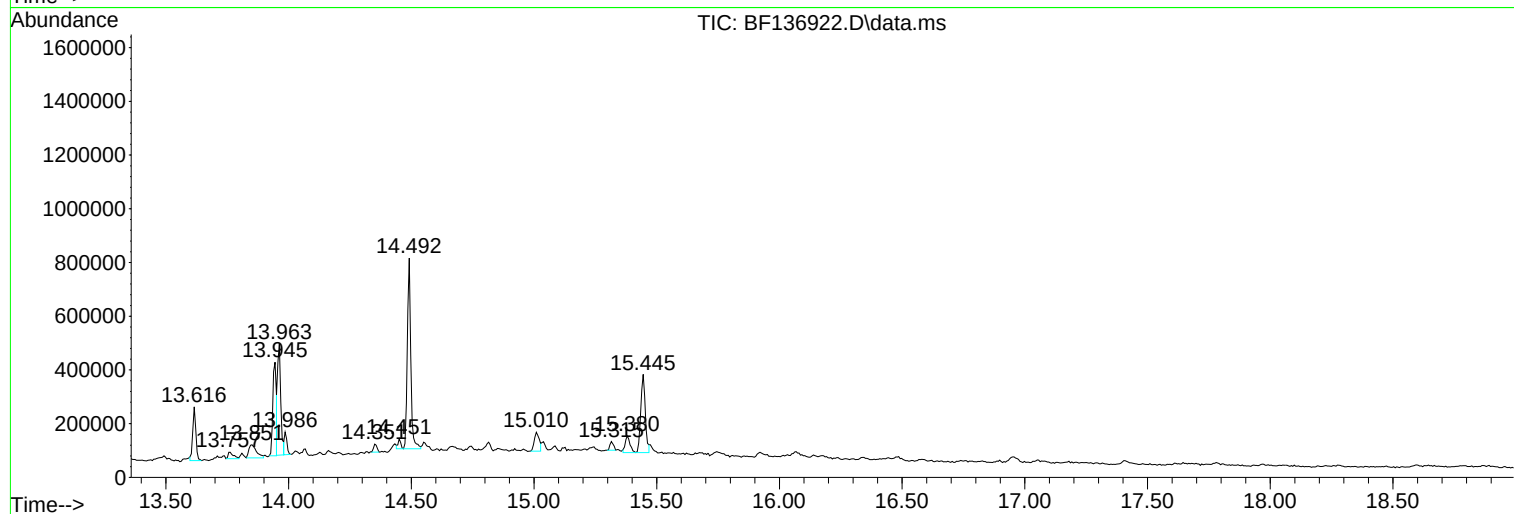
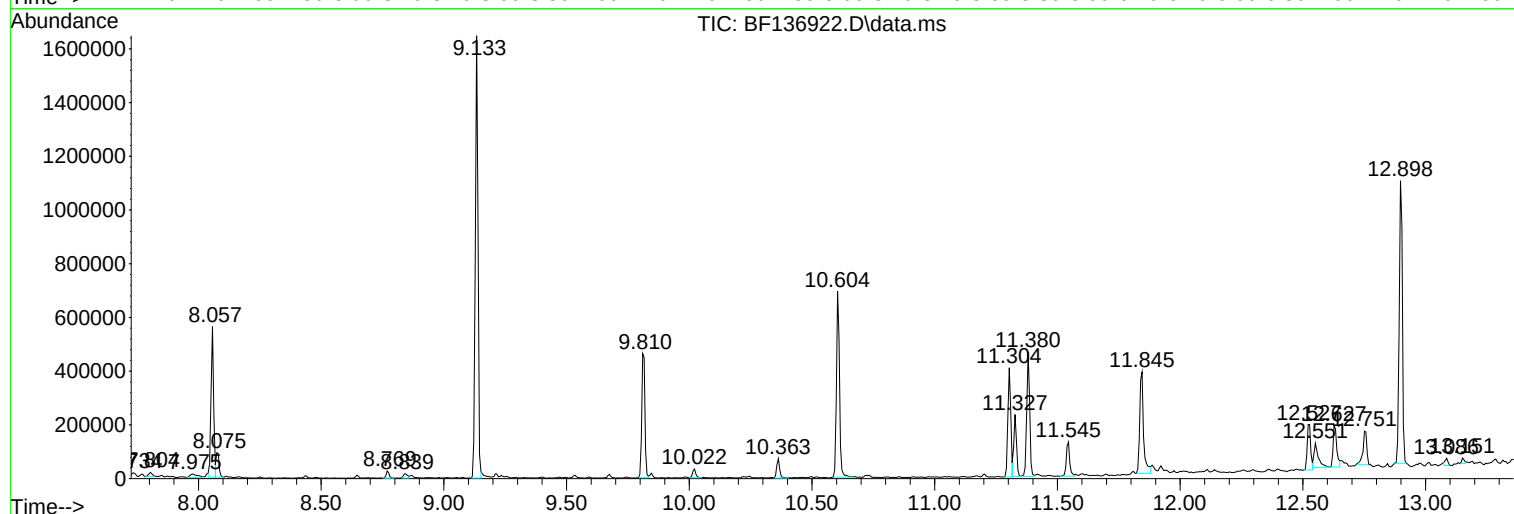
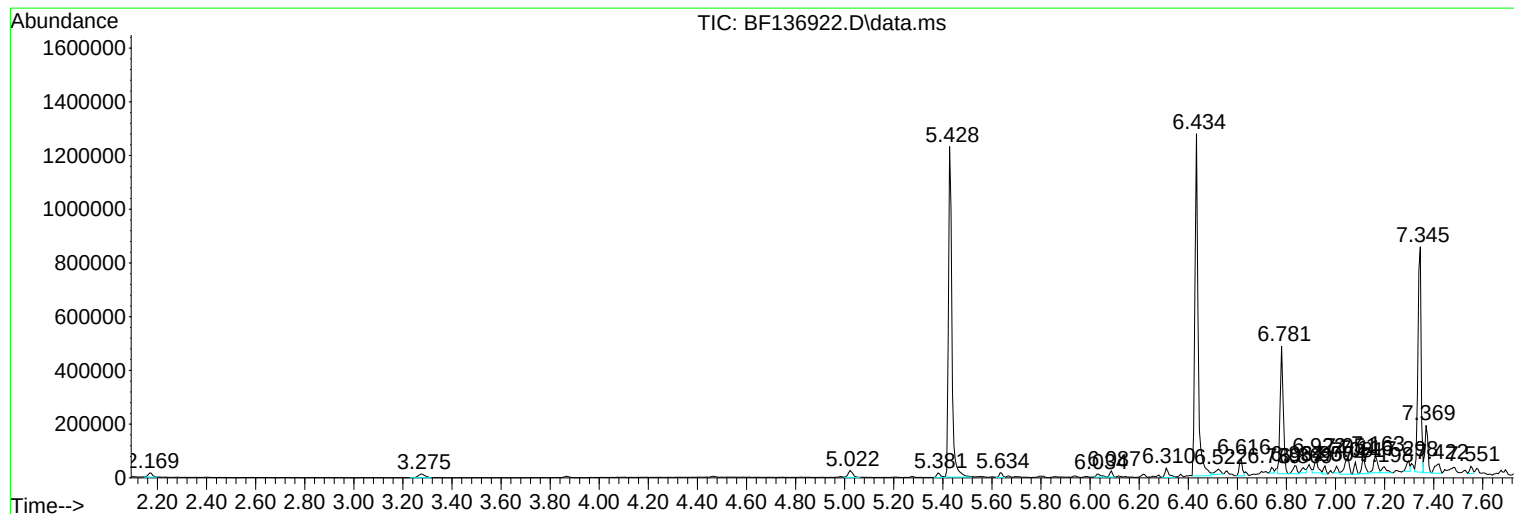
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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Instrument :
BNA_F
ClientSampleId :
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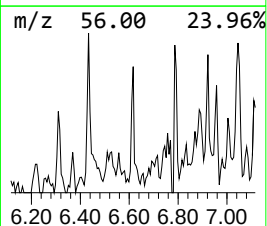
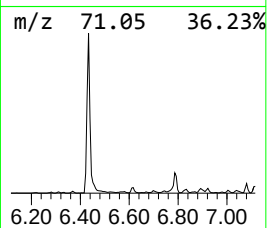
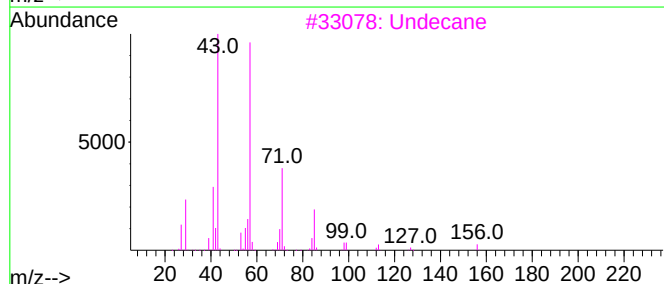
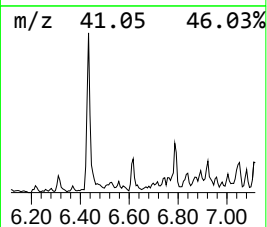
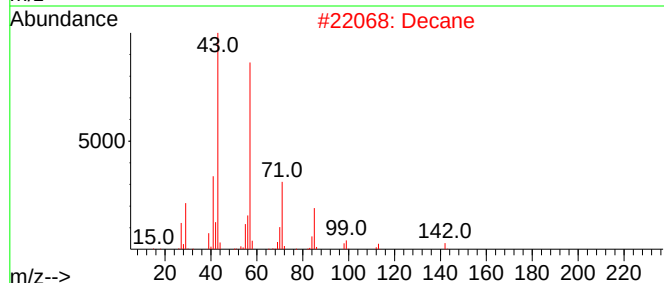
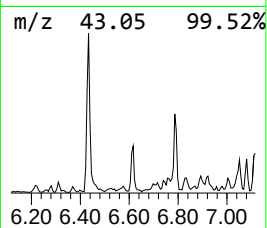
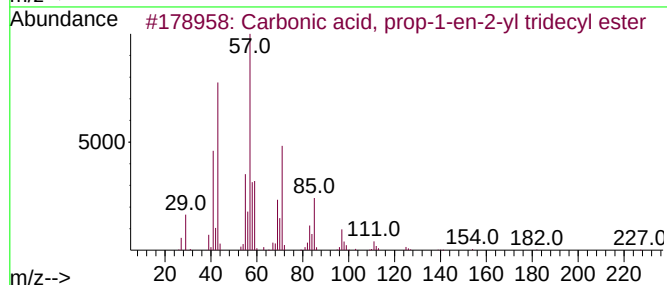
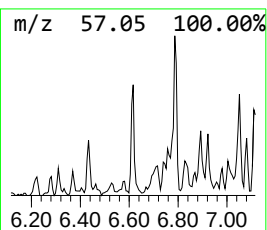
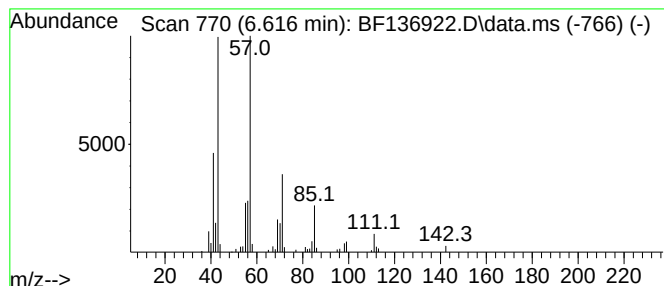
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Carbonic acid, prop-1-en-2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.616	2.70 ng	56872	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	86
2			Decane	142	C10H22	000124-18-5	81
3			Undecane	156	C11H24	001120-21-4	64
4			Carbonic acid, decyl vinyl ester	228	C13H24O3	1000383-25-7	64
5			Nonane, 2-methyl-	142	C10H22	000871-83-0	64



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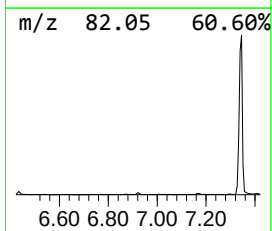
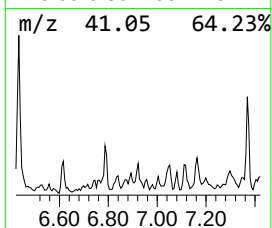
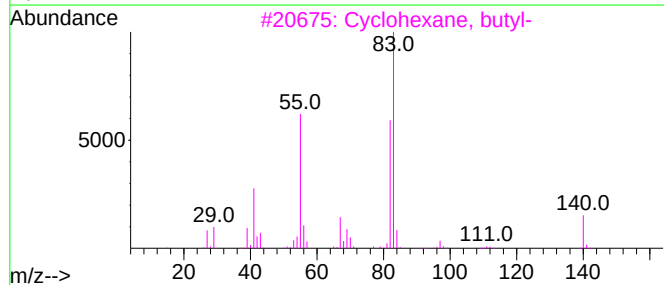
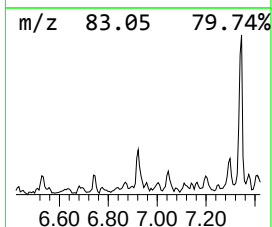
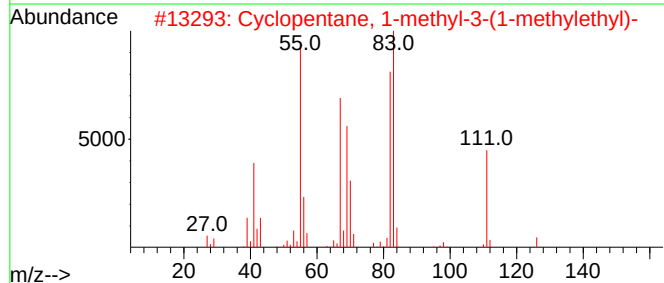
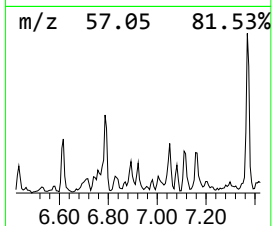
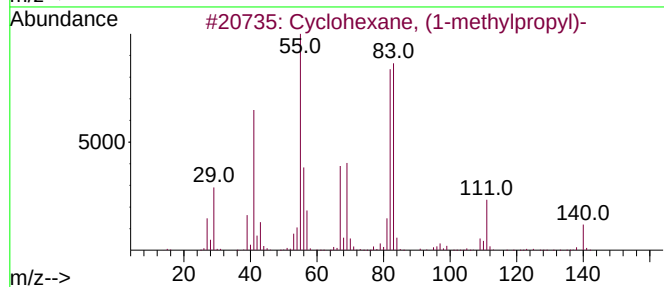
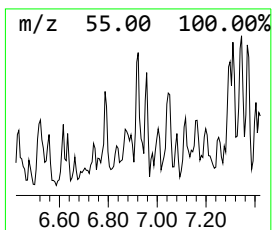
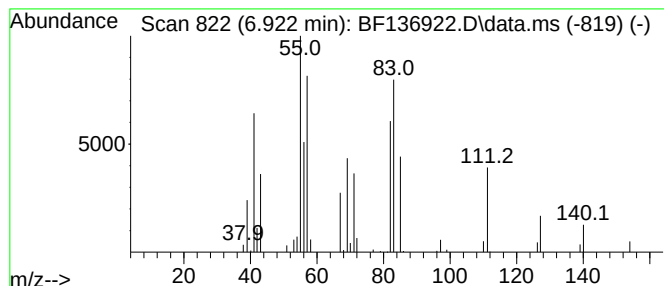
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 2 unknown6.922 Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.922	2.57 ng	54235	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (1-methylpropyl)-	140	C10H20	007058-01-7	49
2			Cyclopentane, 1-methyl-3-(1-meth...	126	C9H18	053771-88-3	43
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	38
4			Succinic acid, 3-methylbut-2-yl ...	270	C15H26O4	1000391-08-2	38
5			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	35



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Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GHL-SS-08

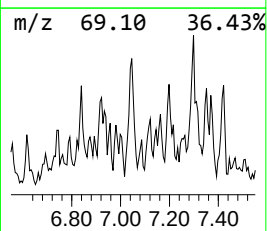
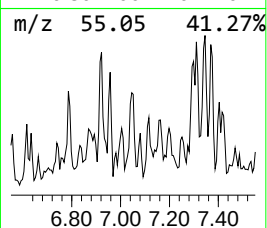
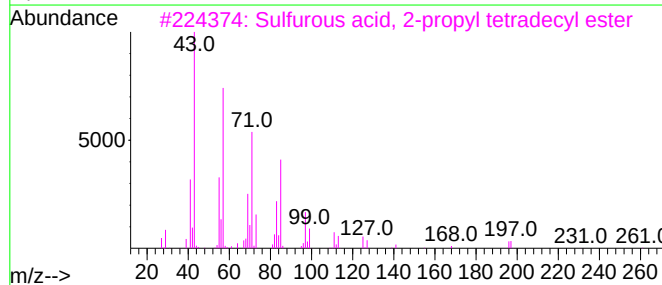
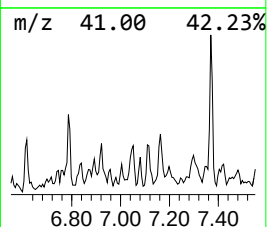
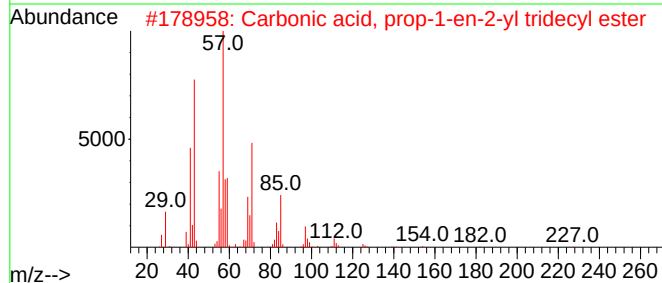
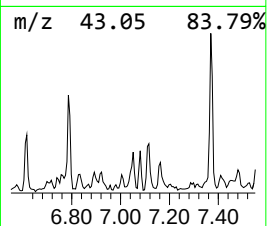
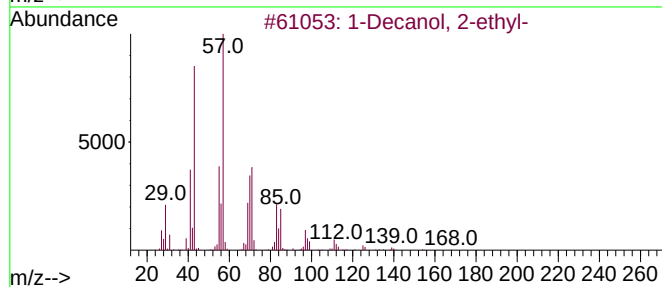
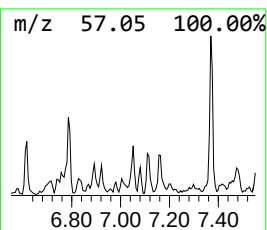
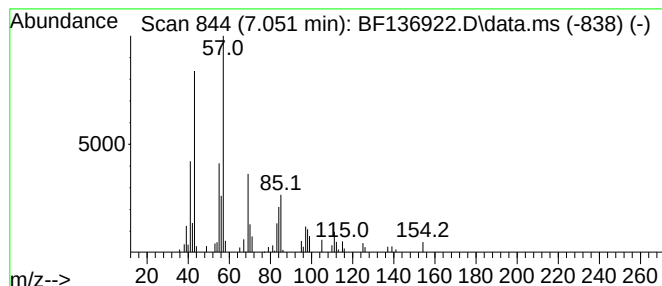
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 1-Decanol, 2-ethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.051	3.87 ng	81579	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Decanol, 2-ethyl-	186	C12H26O	021078-65-9	64
2			Carbonic acid, prop-1-en-2-yl tr...	284	C17H32O3	1000382-90-8	47
3			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	43
4			Sulfurous acid, pentadecyl 2-pro...	334	C18H38O3S	1000309-12-6	38
5			Heptane, 4-ethyl-	128	C9H20	002216-32-2	38



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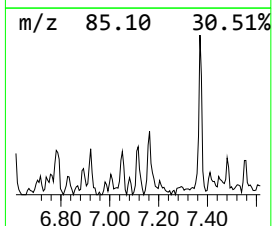
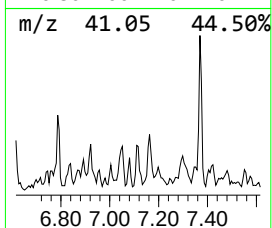
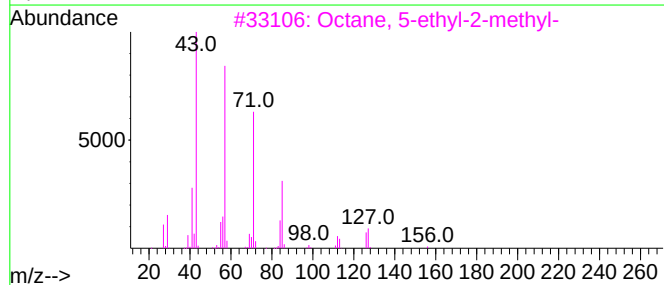
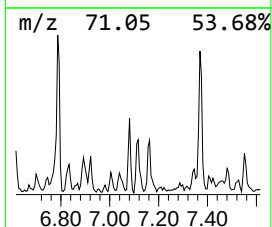
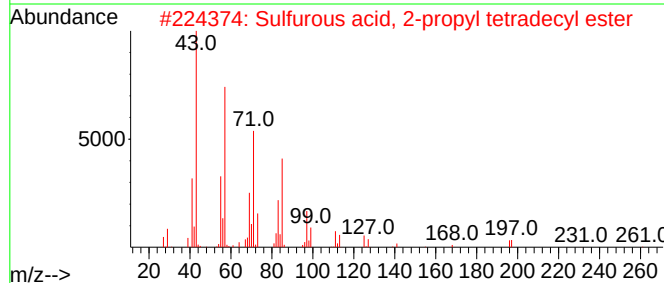
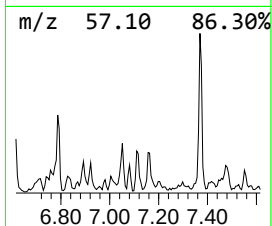
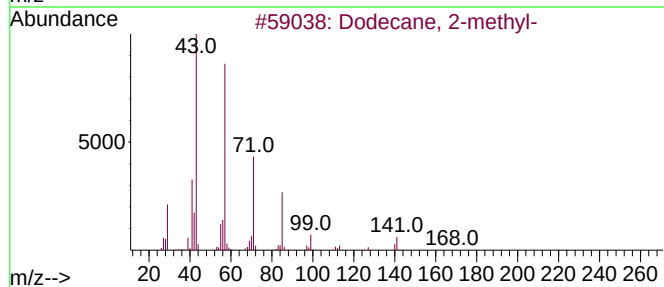
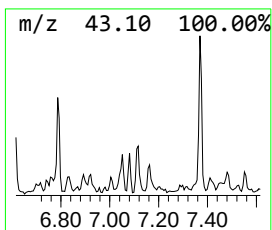
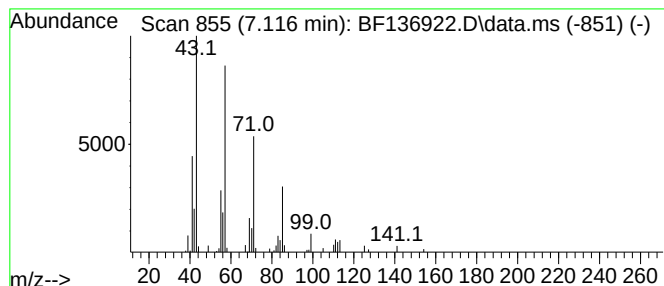
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Dodecane, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.116	2.70 ng	57021	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 2-methyl-	184	C13H28	001560-97-0	72
2			Sulfurous acid, 2-propyl tetradec...	320	C17H36O3S	1010309-12-5	64
3			Octane, 5-ethyl-2-methyl-	156	C11H24	062016-18-6	59
4			Sulfurous acid, dodecyl 2-propyl...	292	C15H32O3S	1000309-12-3	59
5			Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
Operator : CG\JU
Sample : 05899-01
Misc :
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Instrument :
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ClientSampleId :
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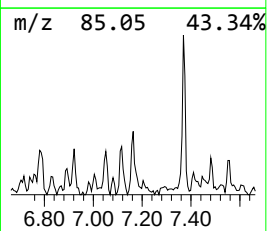
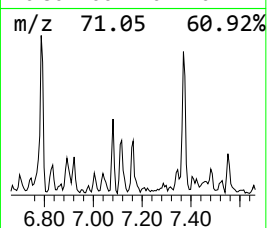
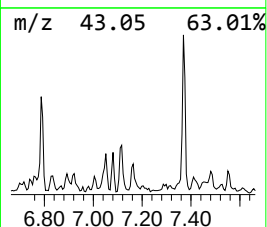
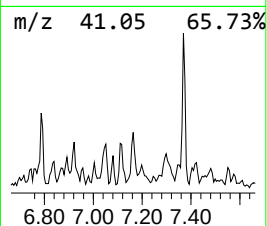
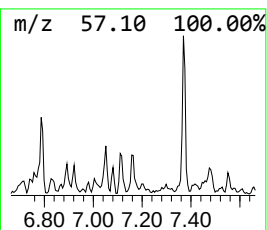
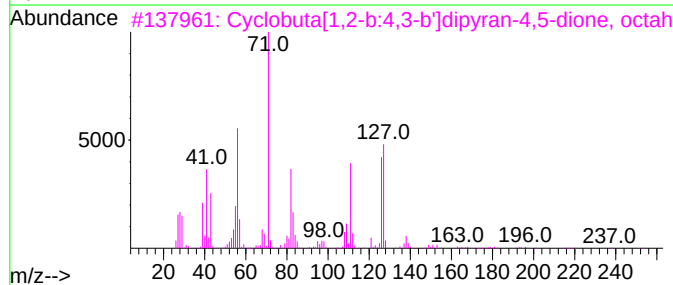
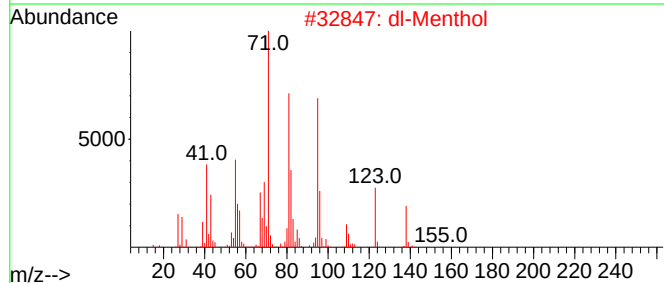
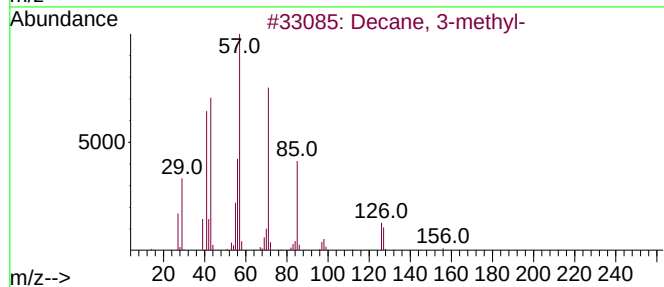
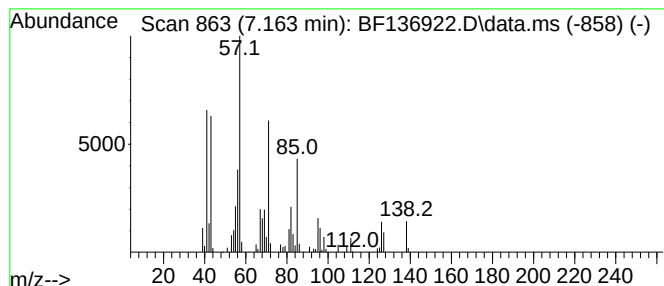
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Decane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.163	3.41 ng	71923	1,4-Dichlorobenzene-d4	6.781

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	50
2			dl-Menthol	156	C10H20O	000089-78-1	43
3			Cyclobuta[1,2-b:4,3-b']dipyrans-4...	252	C14H20O4	054637-09-1	38
4			Cyclohexanol, 5-methyl-2-(1-meth...	156	C10H20O	023283-97-8	38
5			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
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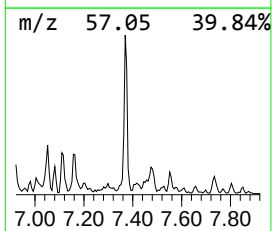
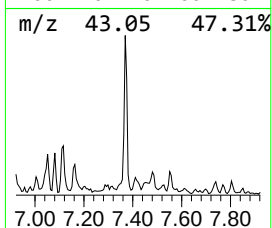
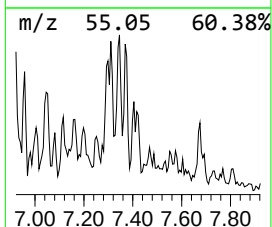
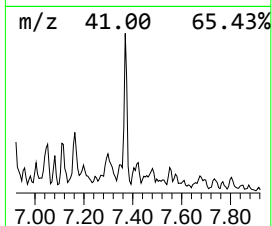
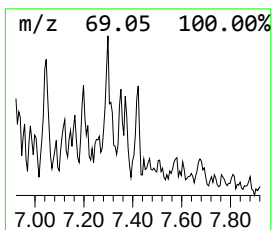
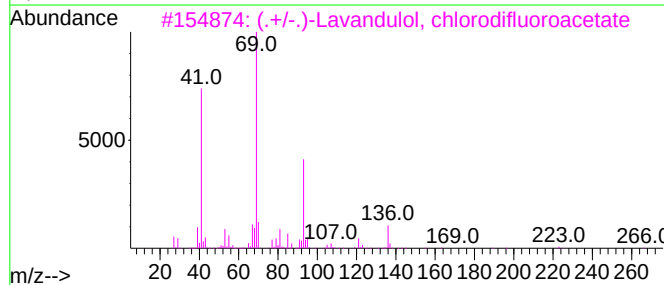
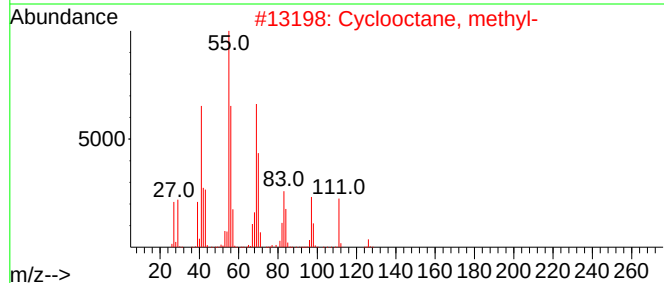
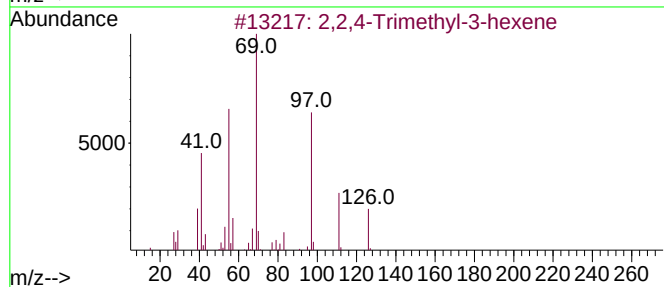
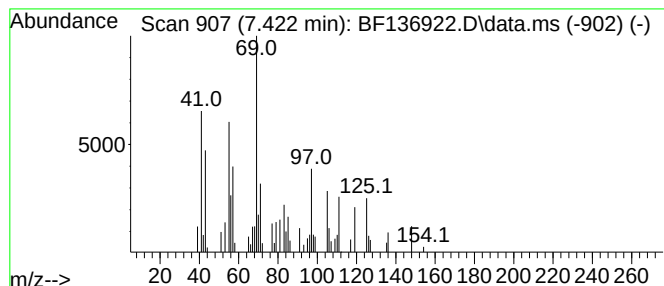
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 unknown7.422 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.422	2.51 ng	52723	Naphthalene-d8	8.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,2,4-Trimethyl-3-hexene	126	C9H18	059643-72-0	46
2			Cyclooctane, methyl-	126	C9H18	001502-38-1	43
3			(.+/-)-Lavandulol, chlorodifluo...	266	C12H17ClF2O2	1000376-22-9	27
4			Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	27
5			Cyclopentane, 1-butyl-2-ethyl-	154	C11H22	072993-32-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
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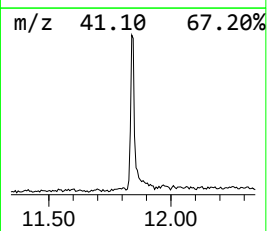
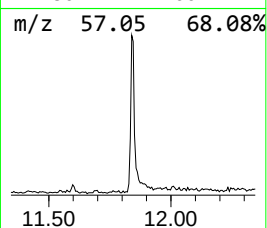
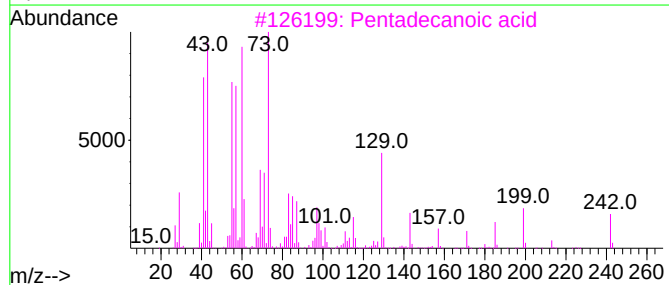
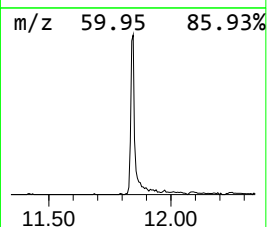
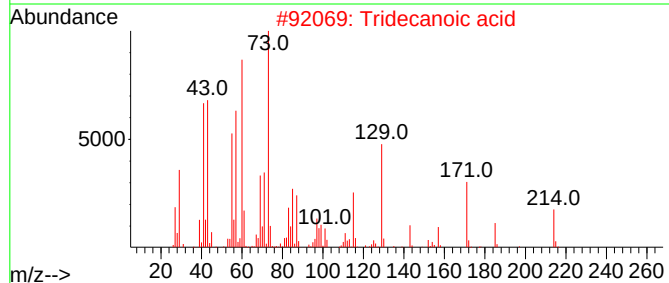
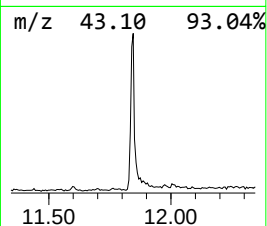
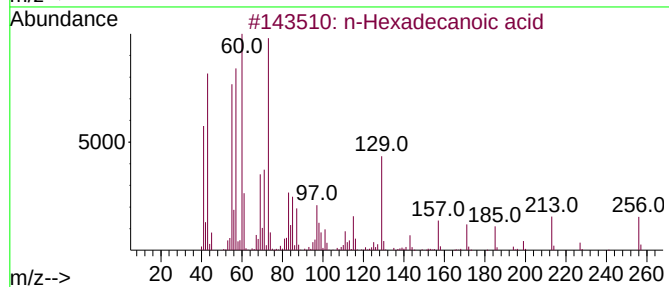
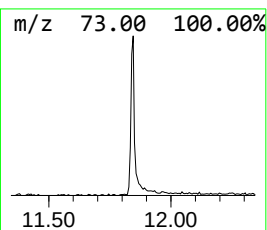
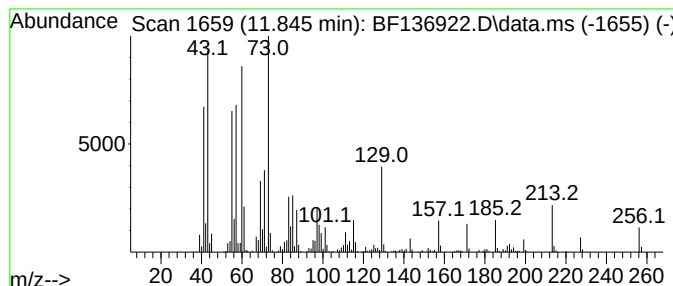
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.845	24.33 ng	413178	Phenanthrene-d10		11.304	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	Tridecanoic acid		214	C13H26O2	000638-53-9	93
3	Pentadecanoic acid		242	C15H30O2	001002-84-2	91
4	Dodecanoic acid		200	C12H24O2	000143-07-7	68
5	Tetradecanoic acid		228	C14H28O2	000544-63-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
Operator : CG\JU
Sample : 05899-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

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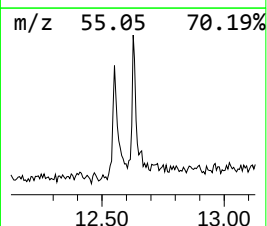
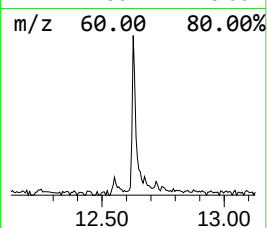
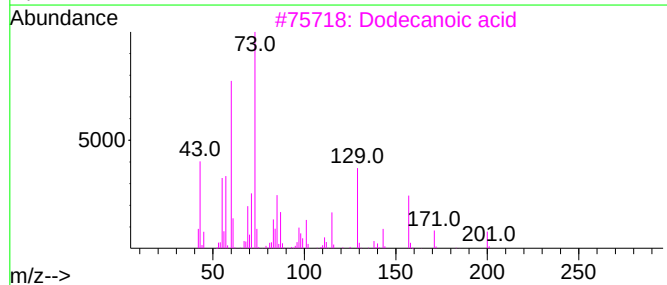
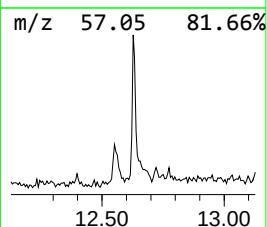
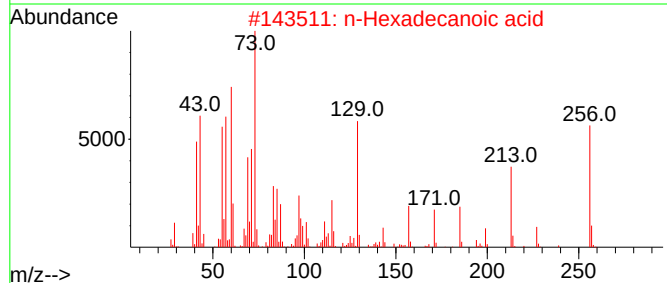
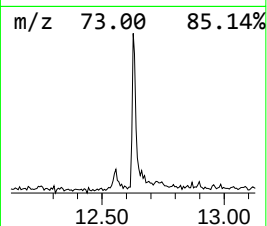
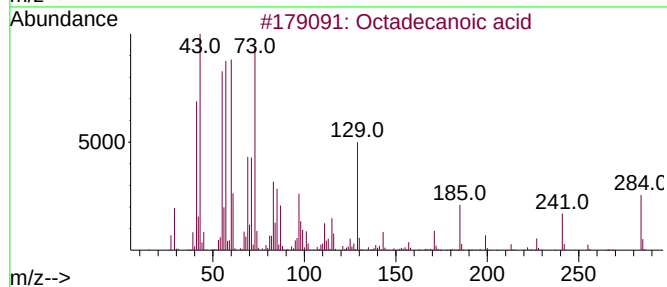
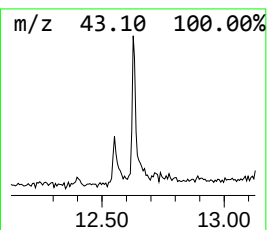
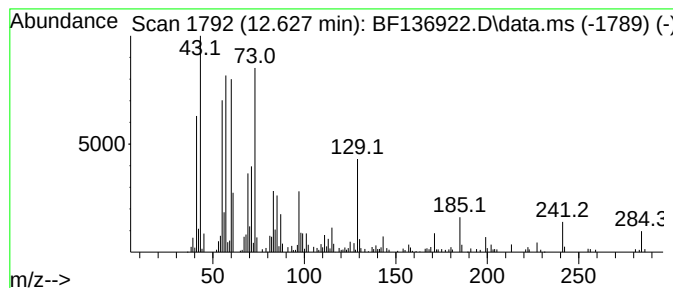
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Octadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.627	9.41 ng	159864	Phenanthrene-d10	11.304

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
3			Dodecanoic acid	200	C12H24O2	000143-07-7	76
4			Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5			Undecanoic acid	186	C11H22O2	000112-37-8	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
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Misc :
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ClientSampleId :
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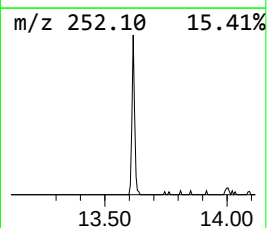
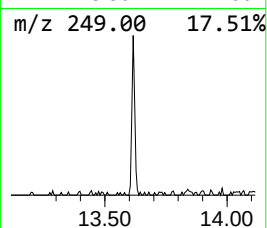
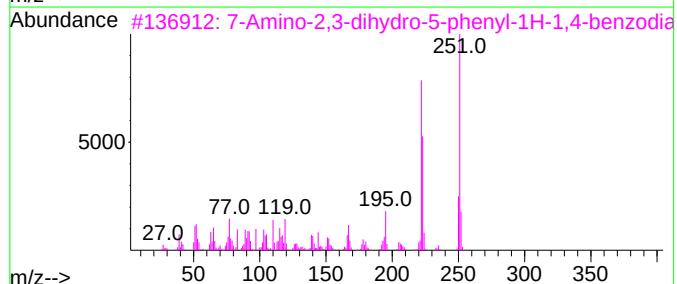
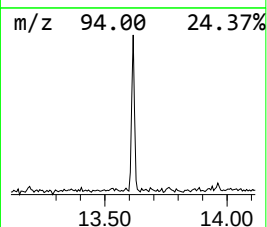
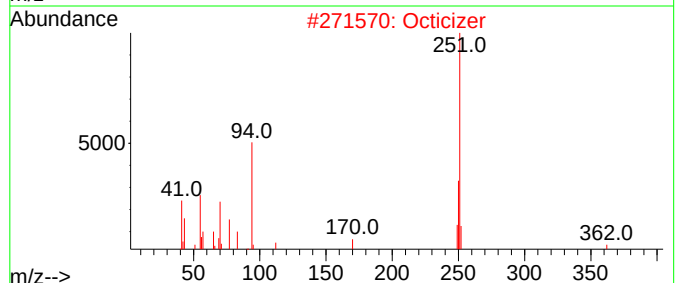
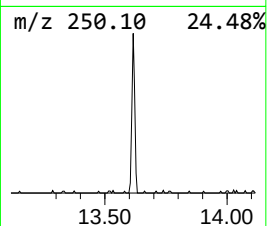
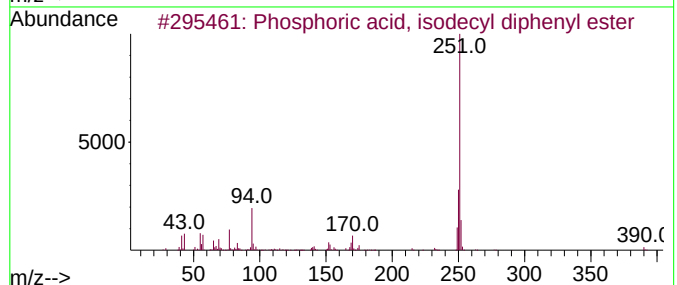
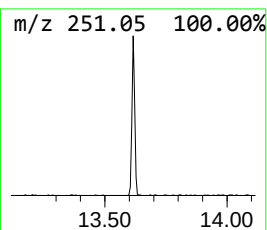
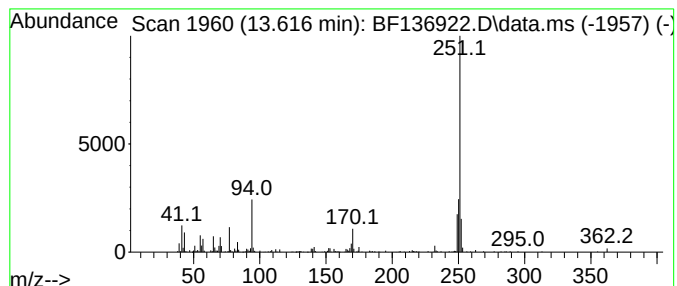
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Phosphoric acid, isodecyl d... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.616	9.47 ng	177620	Chrysene-d12	13.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phosphoric acid, isodecyl diphen...	390	C22H31O4P	1346599-15-2	83
2			Octicizer	362	C20H27O4P	001241-94-7	43
3			7-Amino-2,3-dihydro-5-phenyl-1H-...	251	C15H13N3O	004928-02-3	40
4			2-Benzo[1,3]dioxol-5-yl-7-methyl...	251	C16H13NO2	1000274-75-9	40
5			trans-4'-(Dimethylamino)chalcon	251	C17H17NO	038239-50-8	40



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
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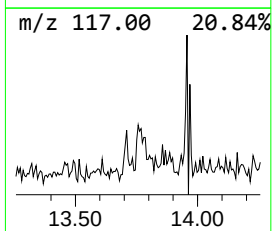
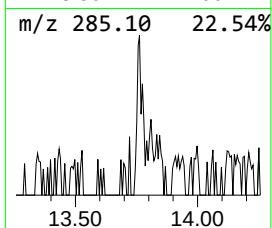
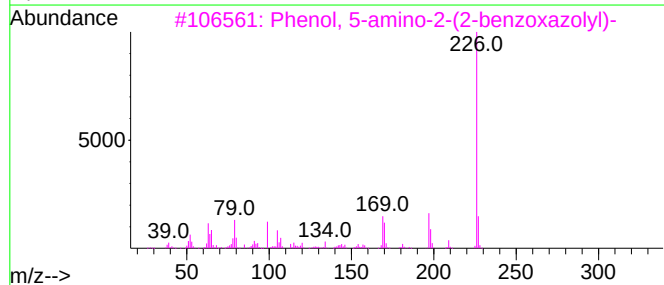
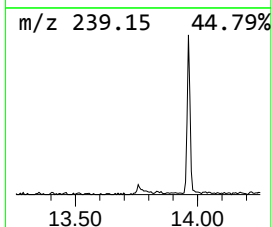
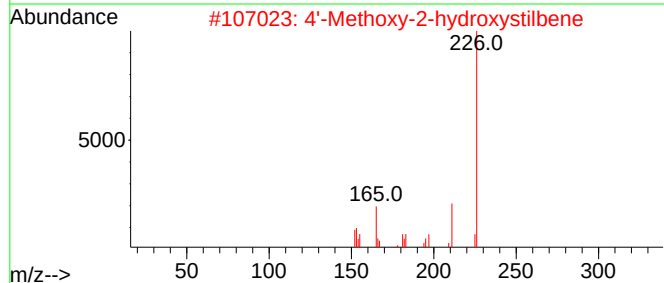
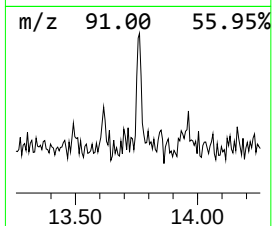
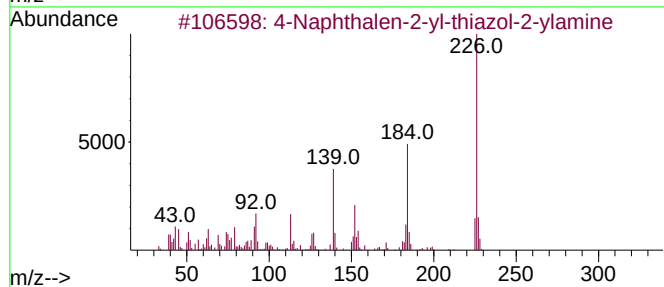
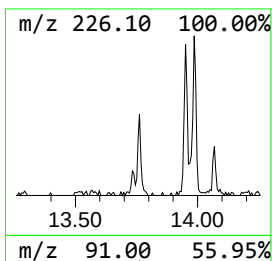
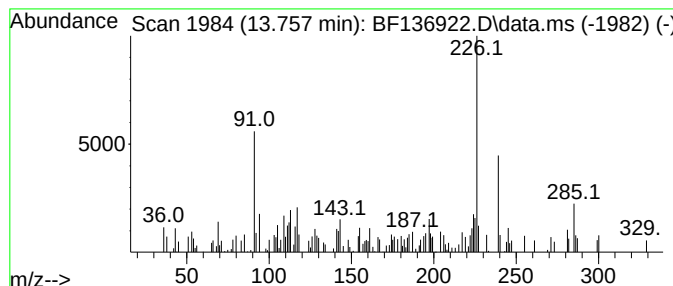
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 unknown13.757 Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.757	2.01 ng	37782	Chrysene-d12	13.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Naphthalen-2-yl-thiazol-2-ylamine	226	C13H10N2S	1000300-53-4	38
2			4'-Methoxy-2-hydroxystilbene	226	C15H14O2	1000147-93-1	38
3			Phenol, 5-amino-2-(2-benzoxazolyl)-	226	C13H10N2O2	088877-61-6	30
4			1,3-Diamino-5,6-dihydro-9-methyl...	226	C13H14N4	037436-39-8	27
5			2,2-Dimethyl-6-hydroxy-7-carbeth...	287	C16H17NO4	031011-05-9	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
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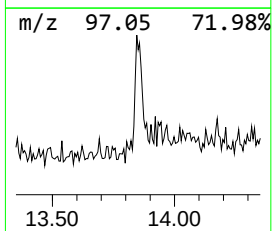
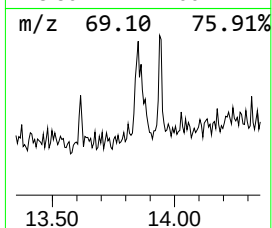
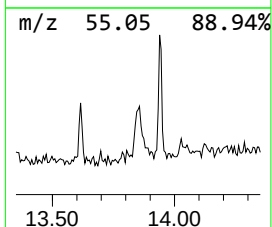
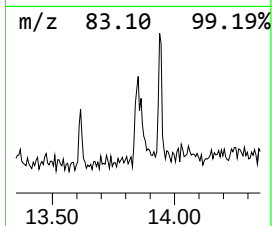
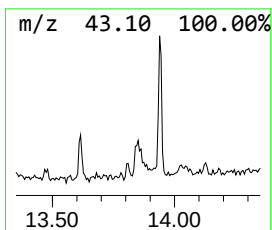
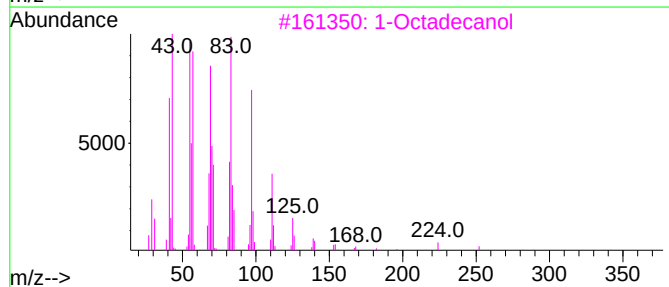
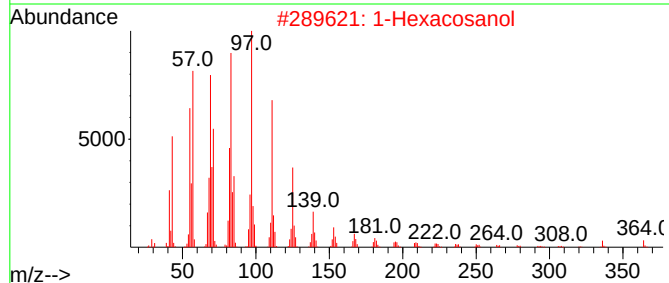
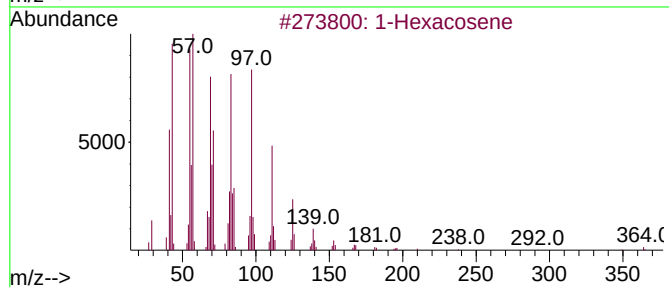
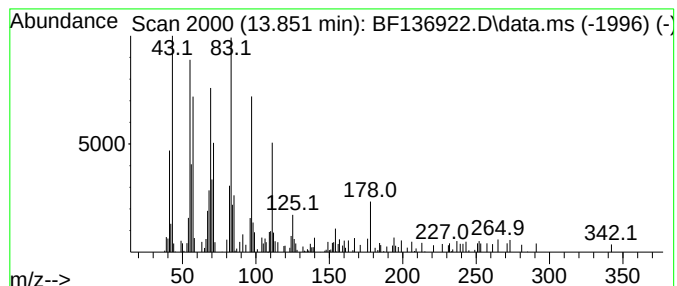
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 1-Hexacosene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.851	5.63 ng	105720	Chrysene-d12	13.963

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexacosene	364	C26H52	018835-33-1	90
2		1-Hexacosanol	382	C26H54O	000506-52-5	89
3		1-Octadecanol	270	C18H38O	000112-92-5	81
4		1-Tetracosene	336	C24H48	010192-32-2	81
5		n-Heptadecanol-1	256	C17H36O	001454-85-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
Operator : CG\JU
Sample : 05899-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GHL-SS-08

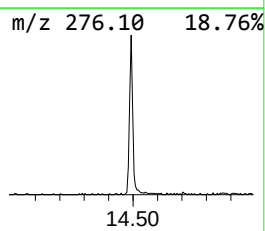
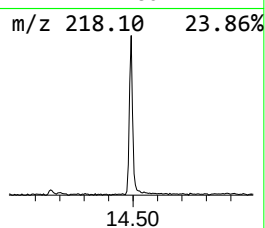
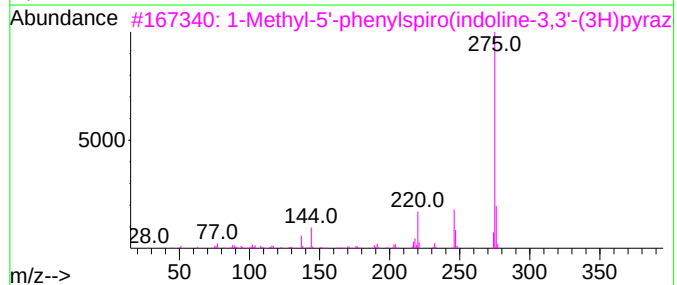
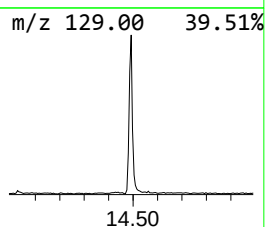
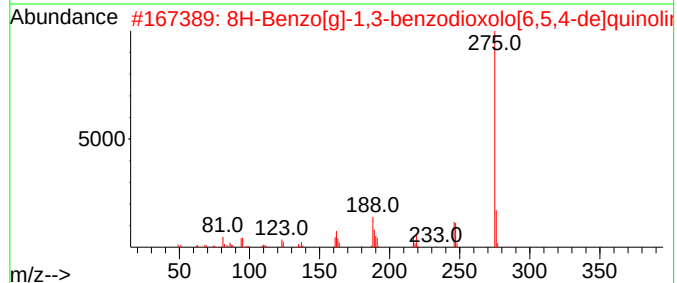
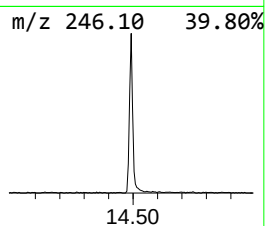
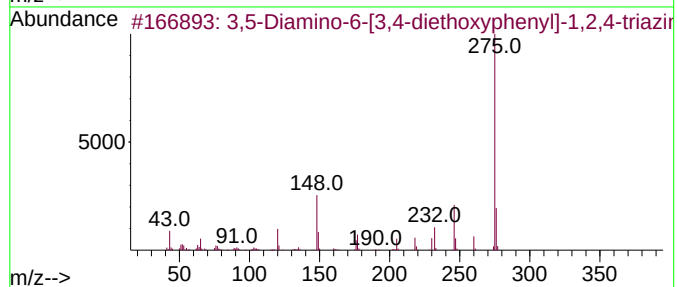
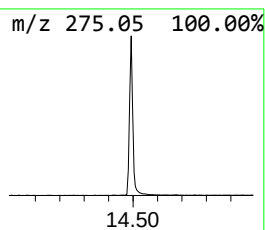
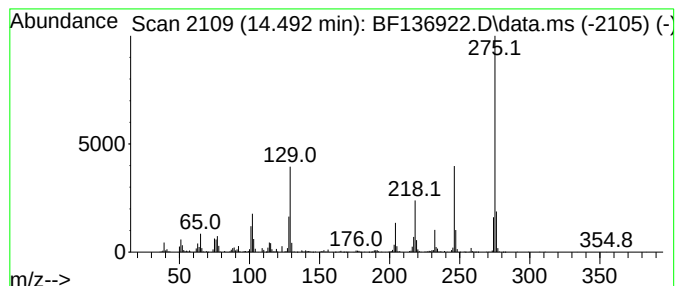
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 unknown14.492 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.492	37.28 ng	699587	Chrysene-d12	13.963

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Diamino-6-[3,4-diethoxypheny...	275	C13H17N5O2	058848-69-4	43
2			8H-Benzo[g]-1,3-benzodioxolo[6,5...	275	C17H9NO3	000475-75-2	38
3			1-Methyl-5'-phenylspiro(indoline...	275	C17H13N3O	015970-21-5	38
4			Acetic acid, 2-[5-(4-nitrophenyl)...	275	C13H13N3O4	351064-45-4	38
5			4H-Isoxazol-5-one, 4-(3,4-dietho...	275	C15H17NO4	1000316-44-0	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF010324\
Data File : BF136922.D
Acq On : 03 Jan 2024 22:57
Operator : CG\JU
Sample : 05899-01
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
GHL-SS-08

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF122023.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
Carbonic acid, ...	6.616	2.7	ng	56872	1	6.781	421856	20.0
unknown6.922	6.922	2.6	ng	54235	1	6.781	421856	20.0
1-Decanol, 2-et...	7.051	3.9	ng	81579	1	6.781	421856	20.0
Dodecane, 2-met...	7.116	2.7	ng	57021	1	6.781	421856	20.0
Decane, 3-methyl-	7.163	3.4	ng	71923	1	6.781	421856	20.0
unknown7.422	7.422	2.5	ng	52723	2	8.057	420796	20.0
n-Hexadecanoic ...	11.845	24.3	ng	413178	4	11.304	339646	20.0
Octadecanoic acid	12.627	9.4	ng	159864	4	11.304	339646	20.0
Phosphoric acid...	13.616	9.5	ng	177620	5	13.963	375303	20.0
unknown13.757	13.757	2.0	ng	37782	5	13.963	375303	20.0
1-Hexacosene	13.851	5.6	ng	105720	5	13.963	375303	20.0
unknown14.492	14.492	37.3	ng	699587	5	13.963	375303	20.0