

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
 Data File : BF137853.D
 Acq On : 08 May 2024 14:12
 Operator : CG\JU
 Sample : P2413-01
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-CC

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-BF043024.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF137853.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.199	16	19	27	rVB2	54754	102819	3.21%	0.371%
2	2.428	51	58	64	rVB	1782301	2370715	74.07%	8.560%
3	3.405	218	224	231	rVB2	31413	67911	2.12%	0.245%
4	3.693	270	273	278	rBV	29276	48481	1.51%	0.175%
5	4.193	352	358	361	rBV2	50905	80324	2.51%	0.290%
6	4.322	376	380	386	rVB	82872	119600	3.74%	0.432%
7	4.463	397	404	408	rBV3	49701	69123	2.16%	0.250%
8	4.599	419	427	430	rBV	29551	42407	1.33%	0.153%
9	4.863	467	472	476	rBV2	59693	79852	2.50%	0.288%
10	5.222	526	533	536	rBV	69857	94195	2.94%	0.340%
11	5.299	542	546	550	rBV3	28367	35680	1.11%	0.129%
12	5.563	587	591	593	rBV2	100903	106456	3.33%	0.384%
13	5.587	593	595	598	rVB	76607	61181	1.91%	0.221%
14	5.675	605	610	613	rBV	2397327	2402459	75.07%	8.675%
15	5.834	631	637	643	rVB6	22529	53012	1.66%	0.191%
16	6.228	701	704	706	rBV	68788	66029	2.06%	0.238%
17	6.293	711	715	717	rBV	39395	41137	1.29%	0.149%
18	6.475	742	746	749	rBV2	28812	38772	1.21%	0.140%
19	6.510	749	752	755	rVV	294650	237998	7.44%	0.859%
20	6.540	755	757	760	rVB	50632	37400	1.17%	0.135%
21	6.569	760	762	765	rBV2	60550	50276	1.57%	0.182%
22	6.604	765	768	770	rVB	44555	36337	1.14%	0.131%
23	6.628	770	772	773	rBV	53263	39863	1.25%	0.144%
24	6.657	773	777	781	rVB	2204744	2447269	76.47%	8.837%
25	6.734	786	790	793	rVB	52535	46682	1.46%	0.169%
26	6.869	810	813	819	rBV	179445	147362	4.60%	0.532%
27	7.051	840	844	847	rBV	918171	816251	25.50%	2.947%
28	7.104	849	853	858	rVB2	69338	72002	2.25%	0.260%
29	7.228	871	874	877	rVB	209090	153458	4.79%	0.554%
30	7.293	882	885	887	rBV	109176	98596	3.08%	0.356%
31	7.363	894	897	903	rVB	236625	248979	7.78%	0.899%
32	7.575	928	933	935	rBV	361874	308521	9.64%	1.114%
33	7.610	935	939	942	rVB	1814802	1607540	50.23%	5.804%
34	7.687	947	952	956	rBV3	37510	41180	1.29%	0.149%
35	7.810	970	973	975	rBV	184183	131686	4.11%	0.475%
36	7.840	975	978	985	rVB2	94751	120600	3.77%	0.435%

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37	7.928	989	993	994	rBV	40118	37419	1.17%	0.135%
38	7.975	998	1001	1003	rBV	55364	49175	1.54%	0.178%
39	7.998	1003	1005	1009	rVB	173363	136805	4.27%	0.494%
40	8.063	1014	1016	1019	rBV	267224	209658	6.55%	0.757%
41	8.334	1049	1062	1064	rBV	1476203	1320496	41.26%	4.768%
42	8.351	1064	1065	1067	rVV	180177	124177	3.88%	0.448%
43	8.375	1067	1069	1072	rVB	68820	52862	1.65%	0.191%
44	8.634	1111	1113	1118	rVB	32693	35318	1.10%	0.128%
45	9.045	1180	1183	1186	rVB	129661	92659	2.90%	0.335%
46	9.145	1195	1200	1203	rBV	79952	67752	2.12%	0.245%
47	9.410	1240	1245	1248	rBV	3697168	3200453	100.00%	11.556%
48	10.098	1358	1362	1365	rBV	1480240	1352701	42.27%	4.884%
49	10.181	1371	1376	1382	rVB	56745	75976	2.37%	0.274%
50	10.886	1492	1496	1499	rBV	2375664	1938424	60.57%	6.999%
51	11.592	1612	1616	1619	rBV	1544078	1328823	41.52%	4.798%
52	12.098	1698	1702	1708	rBV	204661	206444	6.45%	0.745%
53	12.886	1831	1836	1842	rBV3	45045	56149	1.75%	0.203%
54	13.180	1881	1886	1889	rBV	3154322	2958980	92.46%	10.684%
55	14.063	2032	2036	2038	rBV	47322	38561	1.20%	0.139%
56	14.098	2038	2042	2058	rVV6	46898	165520	5.17%	0.598%
57	14.239	2062	2066	2079	rVB	950269	846174	26.44%	3.055%
58	14.692	2139	2143	2147	rVB	37755	38809	1.21%	0.140%
59	15.810	2327	2333	2343	rBV	568019	781077	24.41%	2.820%
60	16.445	2429	2441	2445	rBV	16741	58375	1.82%	0.211%

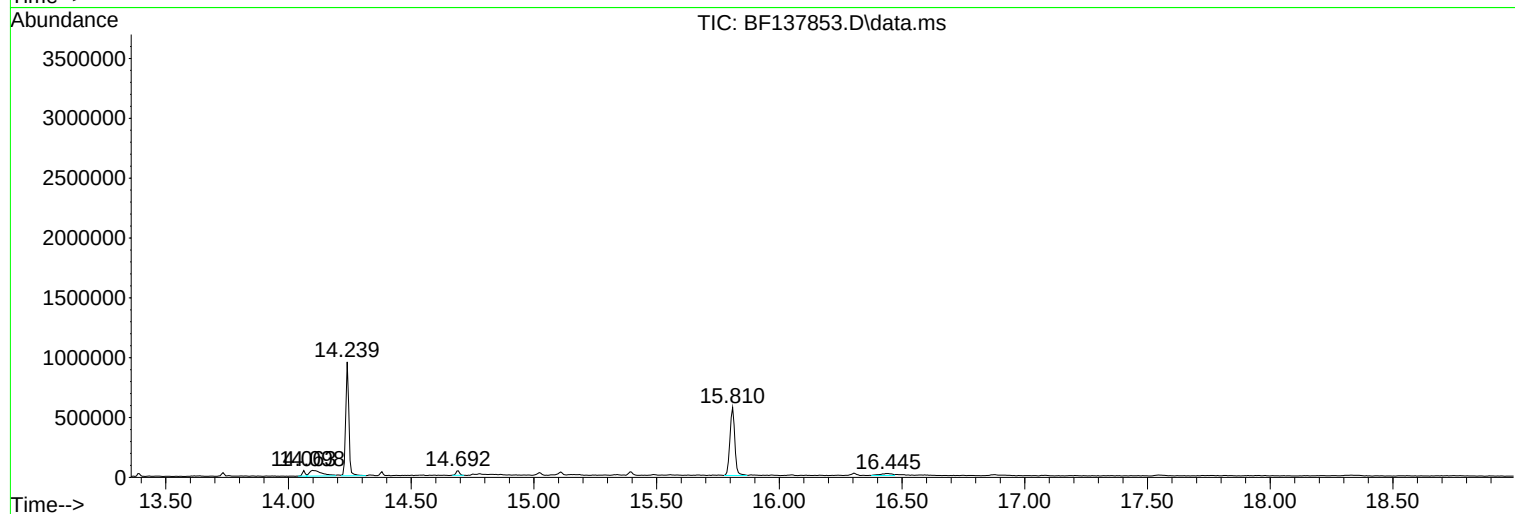
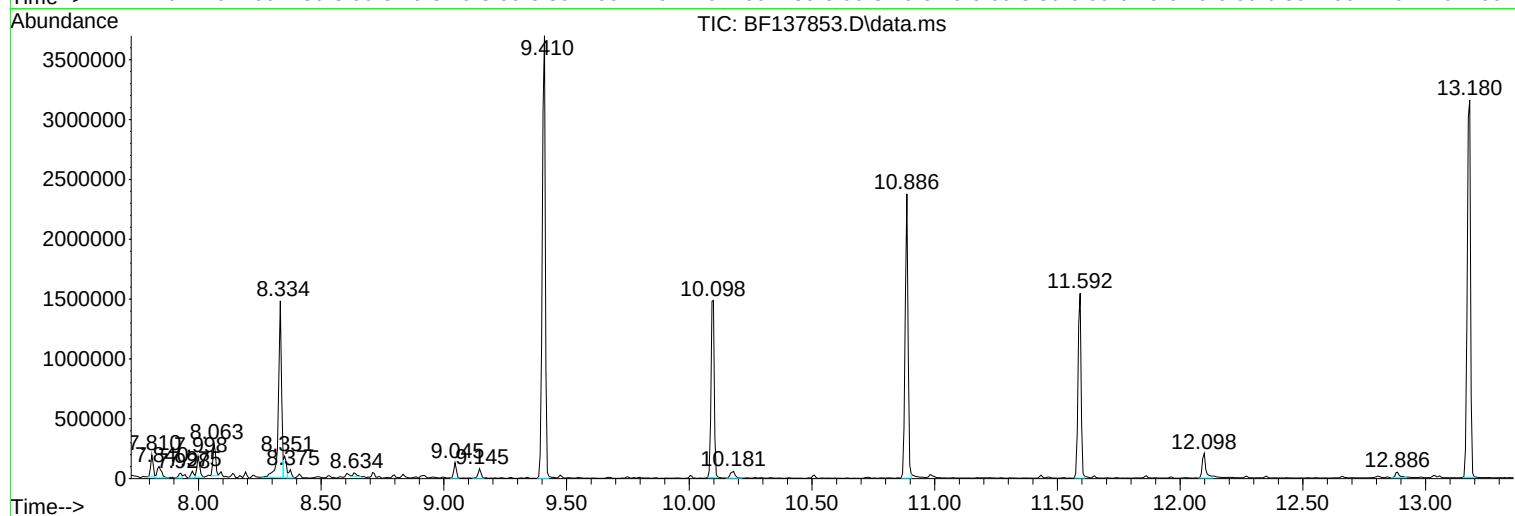
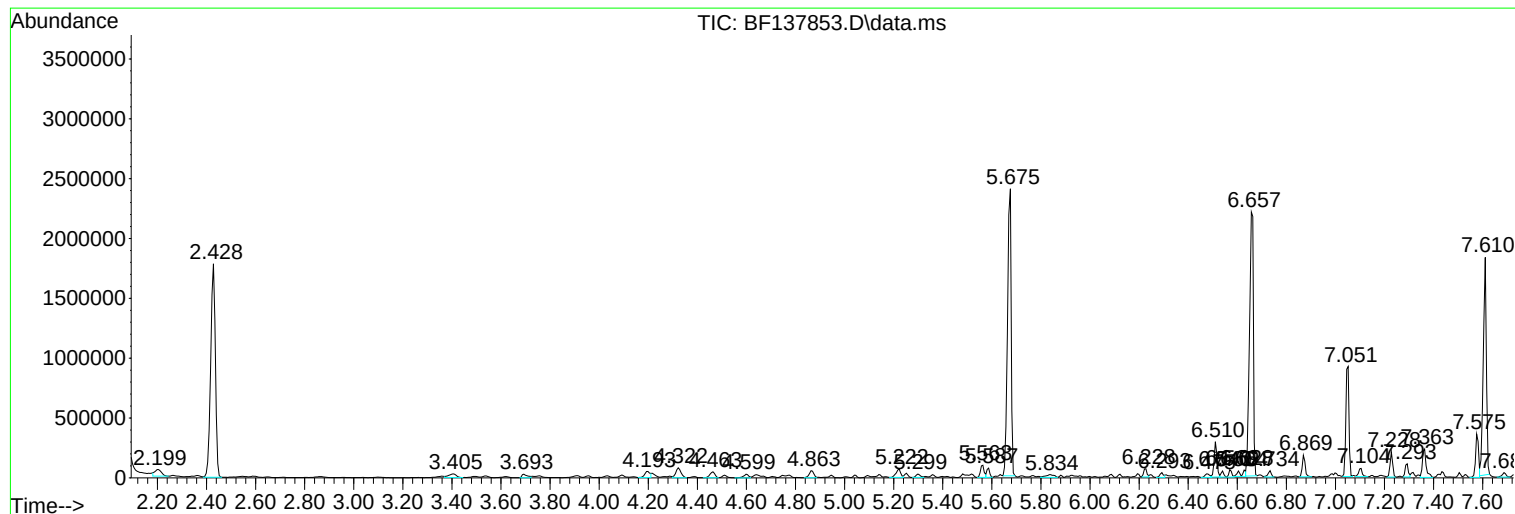
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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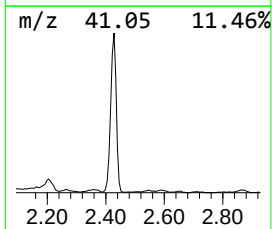
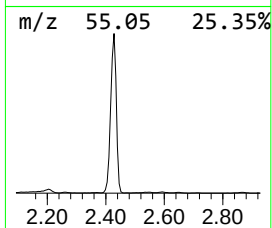
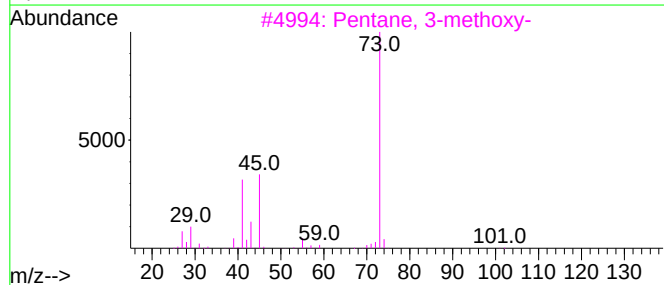
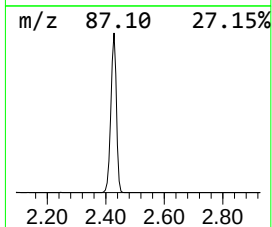
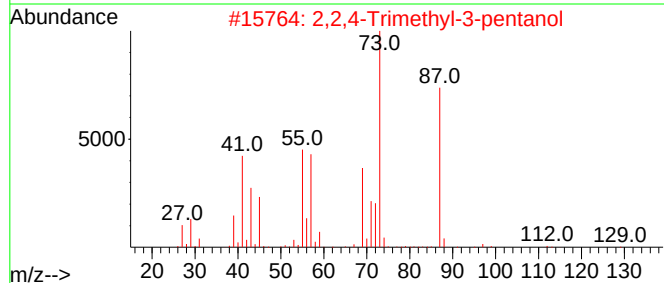
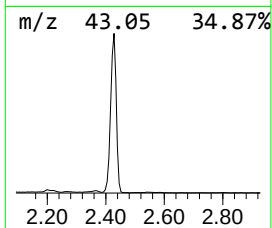
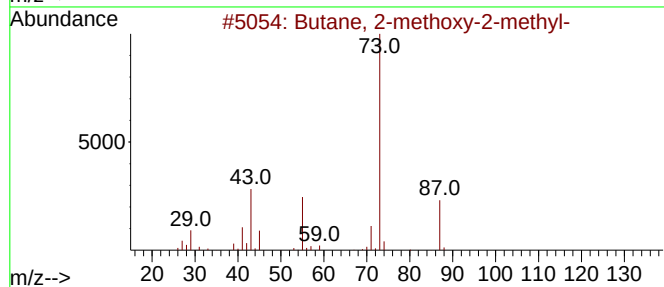
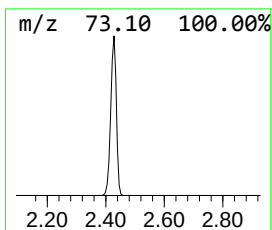
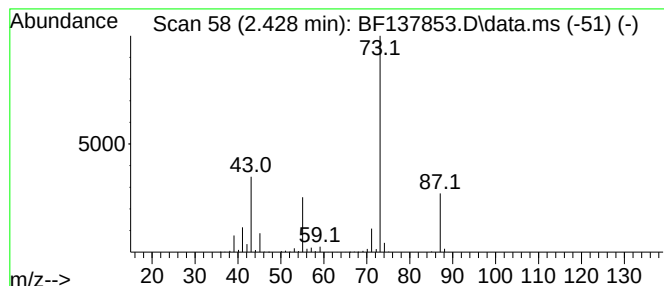
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.428	58.09 ng	2370720	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	78
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Pentane, 3-methoxy-	102	C6H14O	036839-67-5	23
4			Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	9
5			Borane, dimethoxy-	74	C2H7BO2	004542-61-4	9



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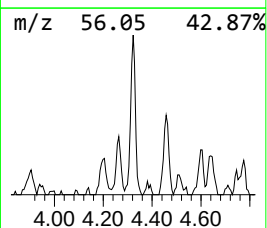
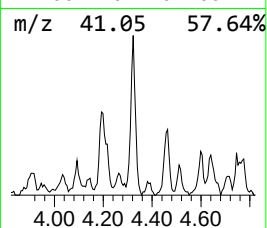
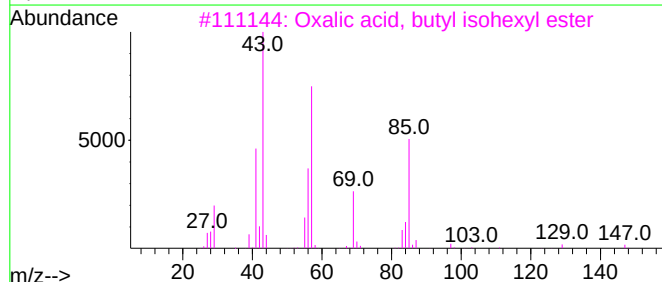
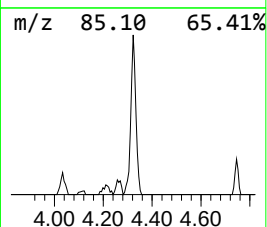
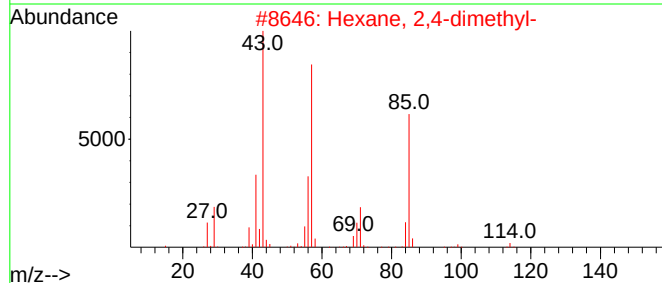
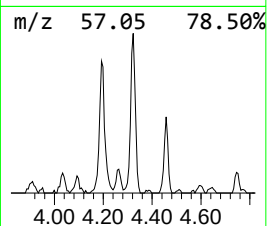
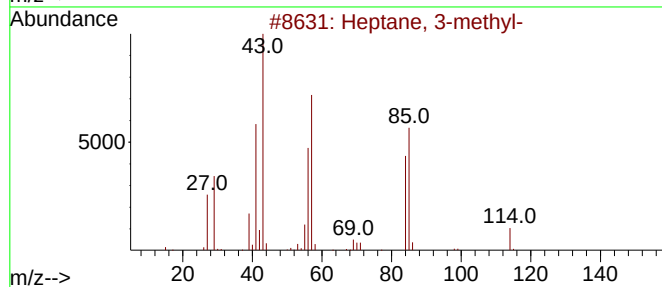
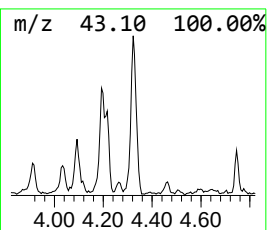
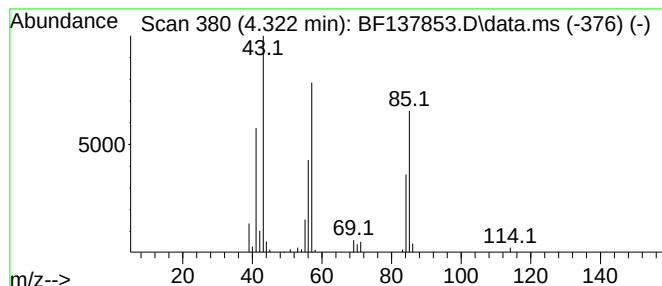
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Heptane, 3-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.322	2.93 ng	119600	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 3-methyl-	114	C8H18	000589-81-1	91
2			Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	80
3			Oxalic acid, butyl isohexyl ester	230	C12H22O4	1000309-32-7	78
4			Hexane, 2-methyl-	100	C7H16	000591-76-4	72
5			Hexane, 1,1'-oxybis-	186	C12H26O	000112-58-3	64



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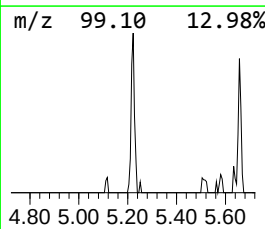
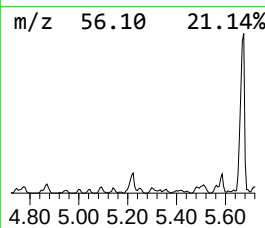
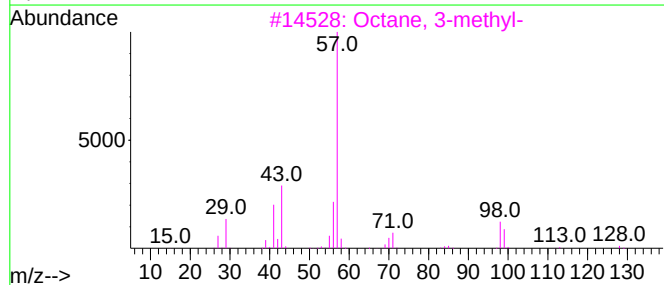
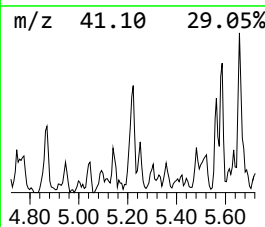
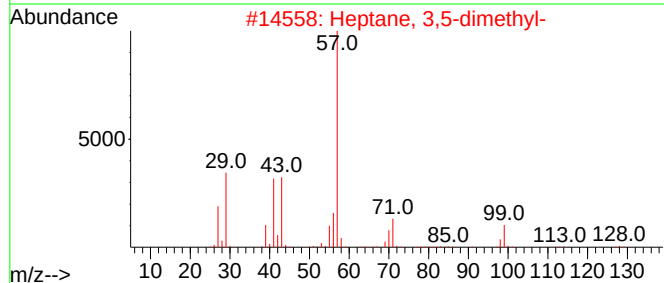
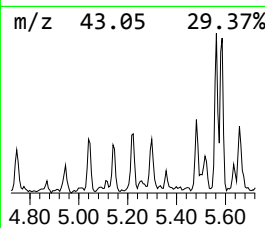
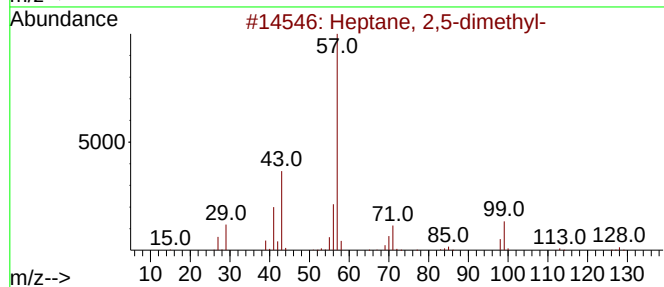
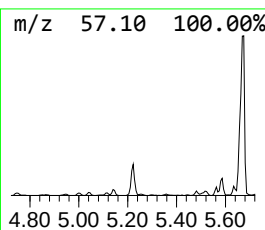
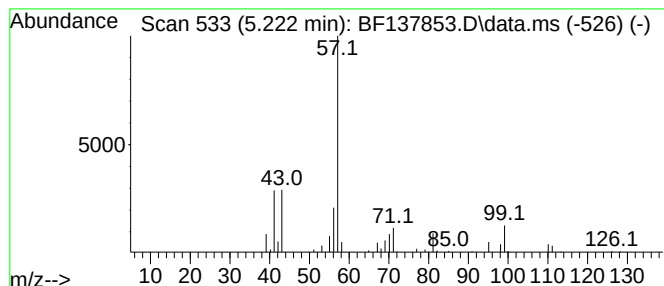
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Heptane, 2,5-dimethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.222	2.31 ng	94195	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptane, 2,5-dimethyl-	128	C9H20	002216-30-0	76
2			Heptane, 3,5-dimethyl-	128	C9H20	000926-82-9	68
3			Octane, 3-methyl-	128	C9H20	002216-33-3	62
4			2,2,4,4-Tetramethyloctane	170	C12H26	062183-79-3	53
5			Decane, 2,6,6-trimethyl-	184	C13H28	062108-24-1	50



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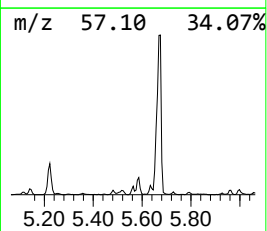
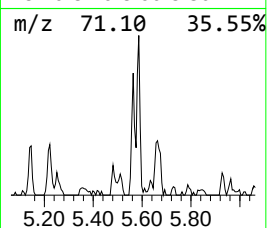
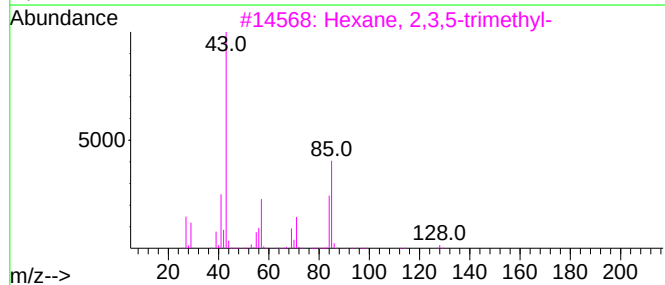
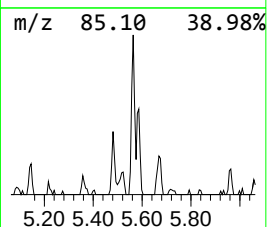
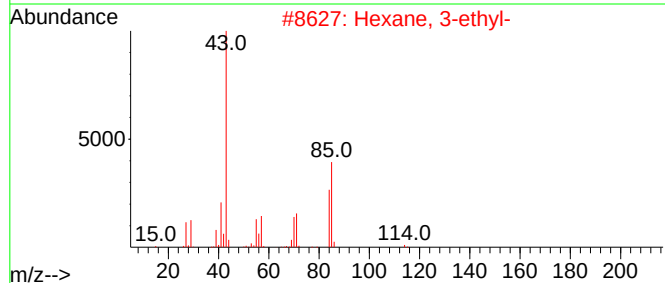
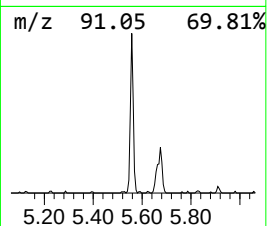
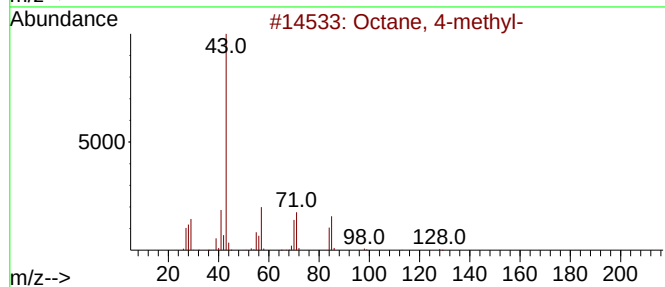
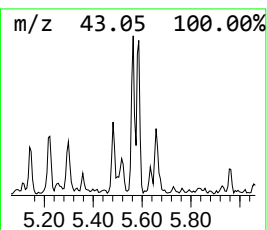
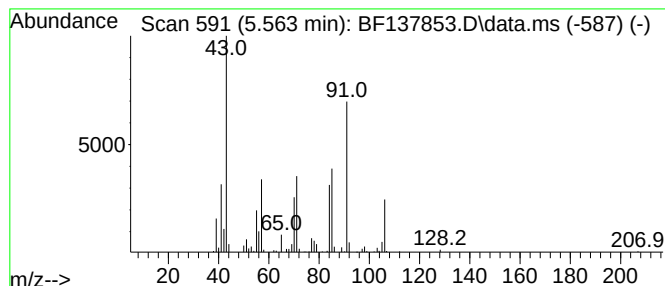
Instrument :
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Octane, 4-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD			R.T.
5.563	2.61 ng	106456	1,4-Dichlorobenzene-d4			7.051
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octane, 4-methyl-		128	C9H20	002216-34-4	81
2	Hexane, 3-ethyl-		114	C8H18	000619-99-8	72
3	Hexane, 2,3,5-trimethyl-		128	C9H20	001069-53-0	52
4	Heptane, 2,4-dimethyl-		128	C9H20	002213-23-2	50
5	Heptane, 4-ethyl-		128	C9H20	002216-32-2	49



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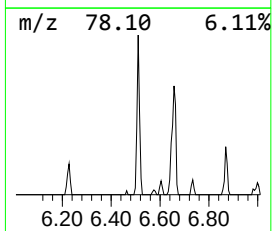
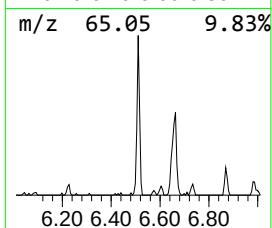
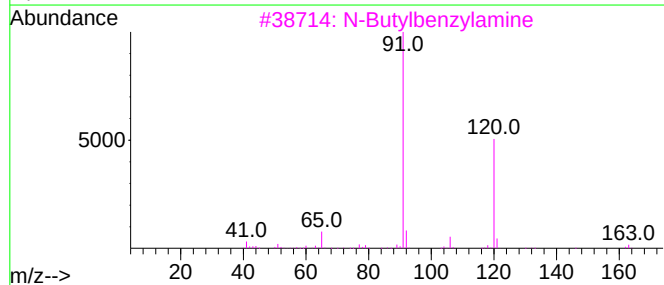
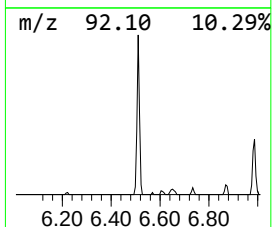
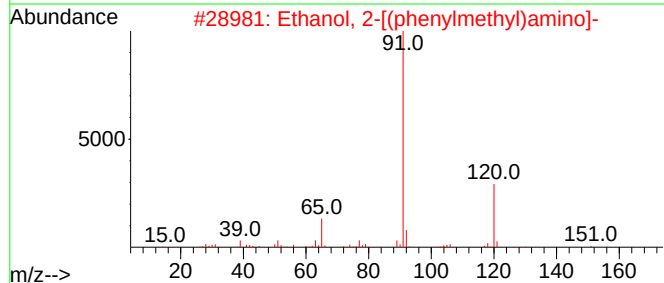
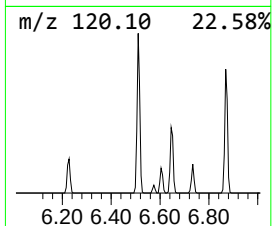
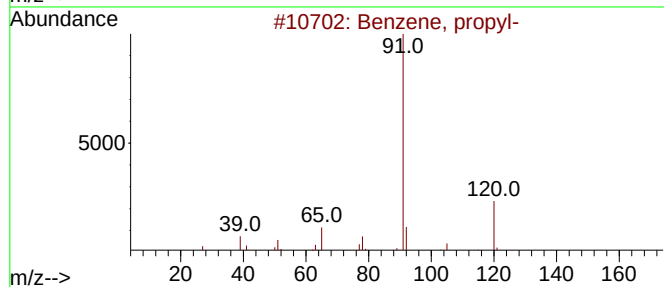
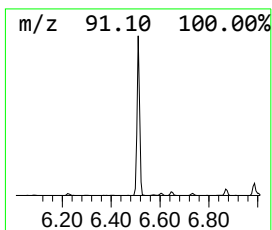
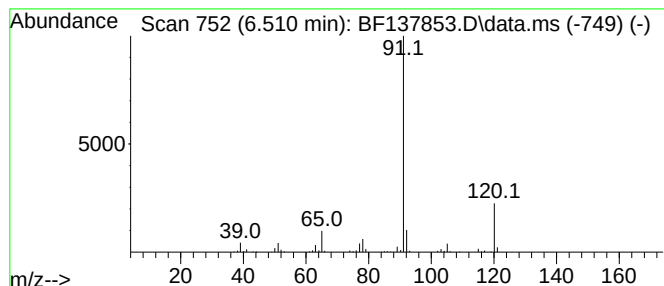
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Benzene, propyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.510	5.83 ng	237998	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Ethanol, 2-[(phenylmethyl)amino]-	151	C9H13NO	000104-63-2	72
3			N-Butylbenzylamine	163	C11H17N	002403-22-7	72
4			Phenethylamine, N-benzyl-p-chloro-	245	C15H16ClN	013622-43-0	72
5			1,2-Ethanediamine, N,N'-bis(phen...	240	C16H20N2	000140-28-3	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-CC

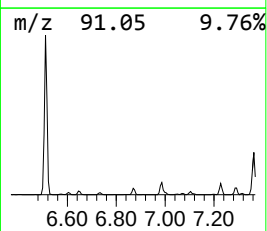
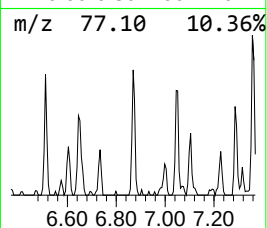
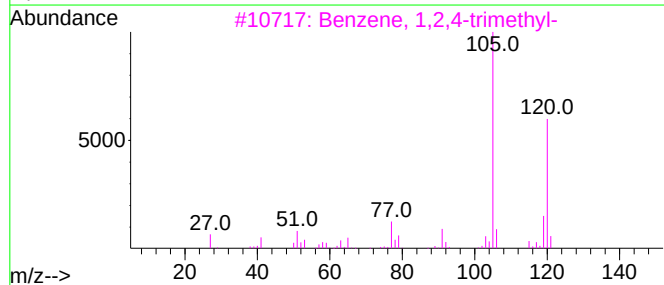
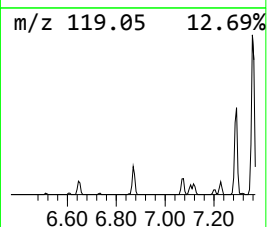
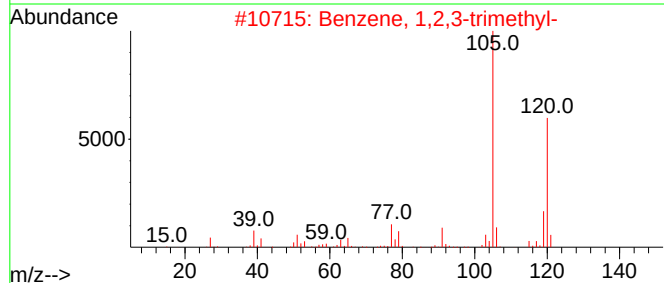
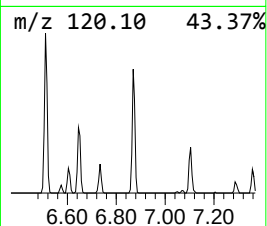
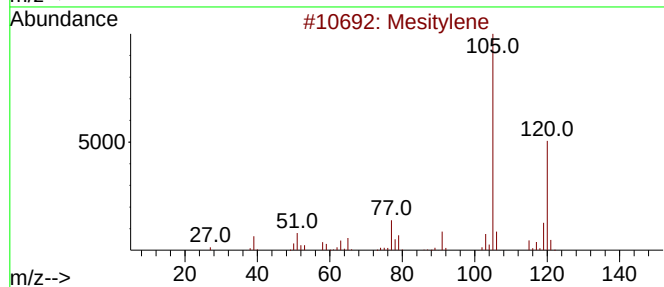
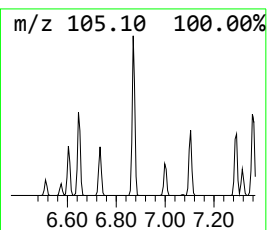
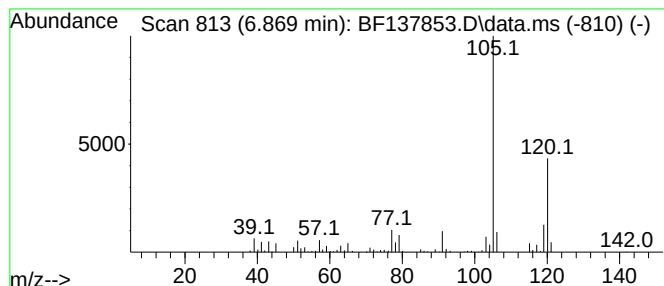
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Mesitylene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.869	3.61 ng	147362	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
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Sample : P2413-01
Misc :
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Instrument :
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ClientSampleId :
WC-CC

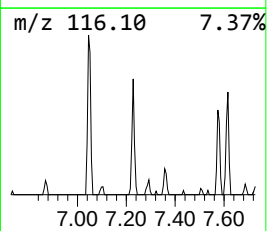
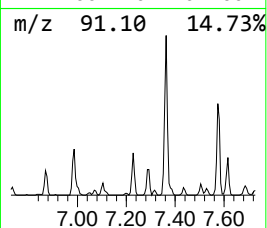
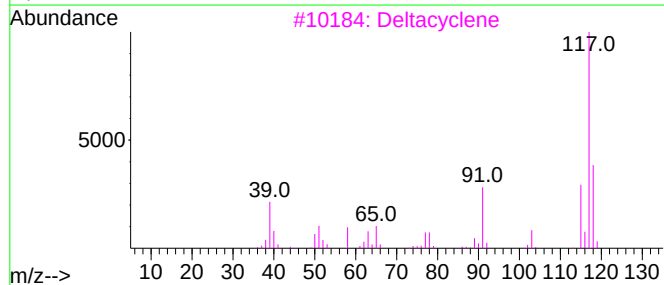
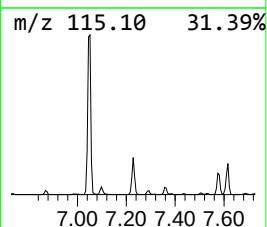
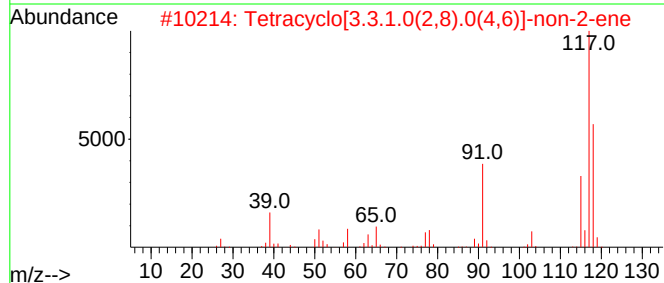
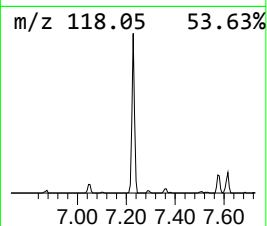
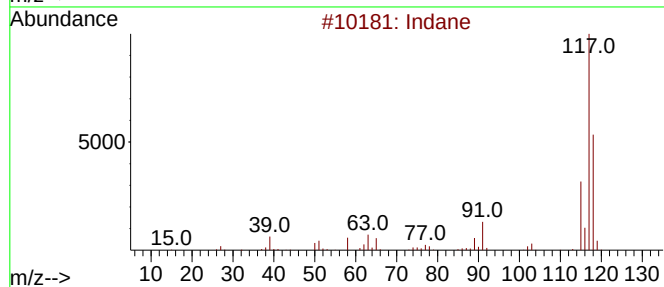
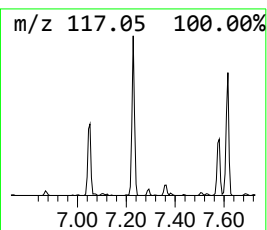
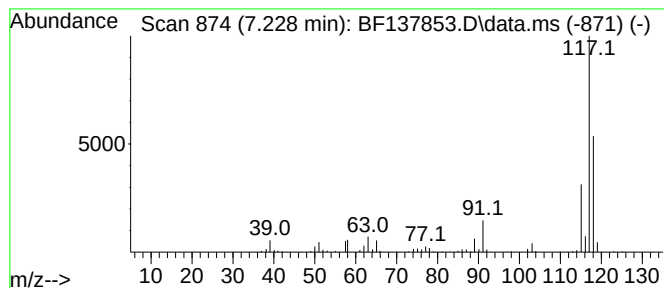
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.228	3.76 ng	153458	1,4-Dichlorobenzene-d4	7.051

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Indane	118	C9H10	000496-11-7	93
2		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	80
3		Deltacyclene	118	C9H10	007785-10-6	72
4		Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5		Benzene, (2-bromocyclopropyl)-	196	C9H9Br	036617-02-4	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-CC

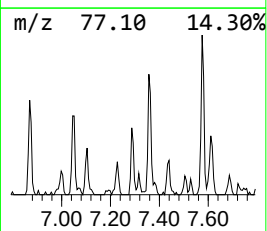
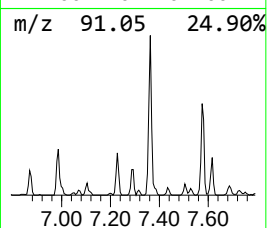
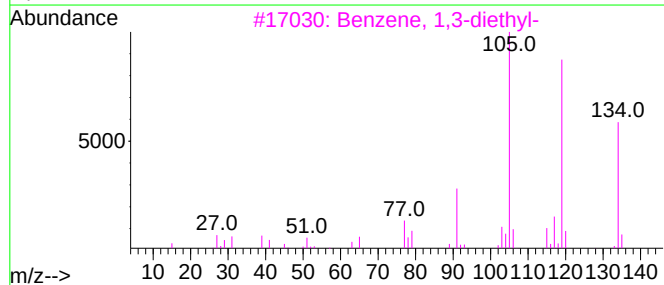
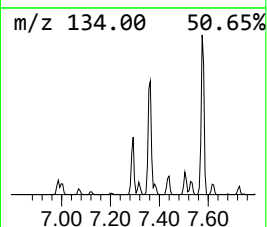
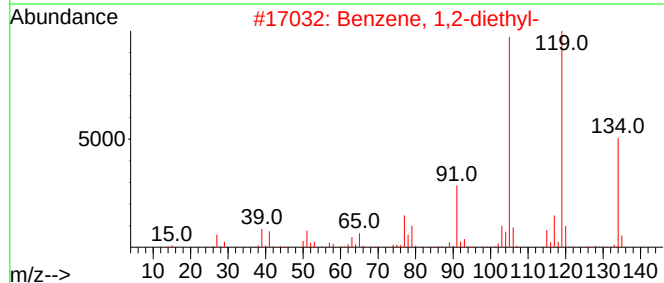
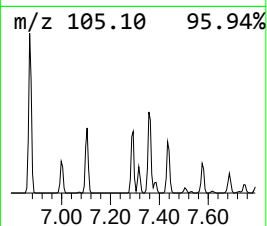
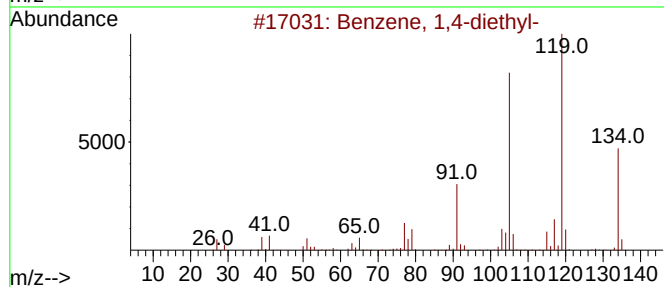
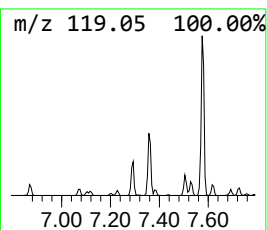
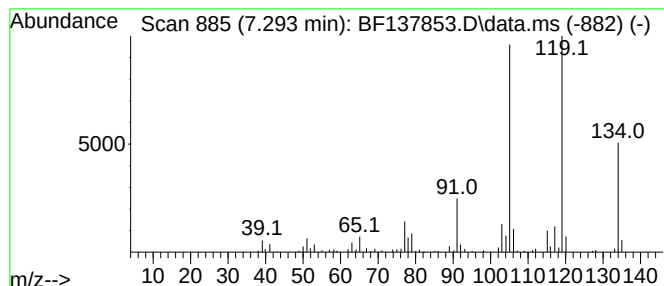
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Benzene, 1,4-diethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.293	2.42 ng	98596	1,4-Dichlorobenzene-d4	7.051

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	90
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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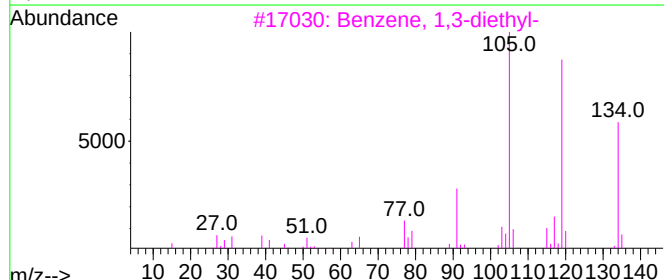
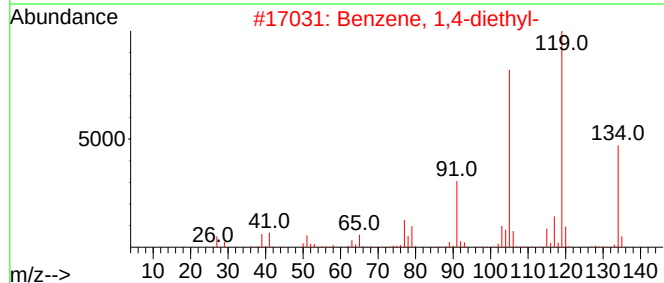
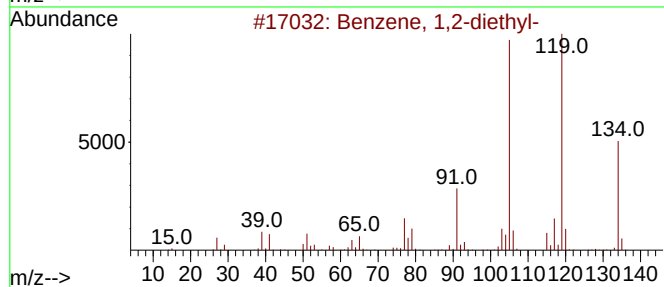
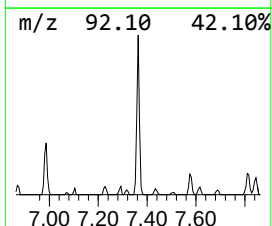
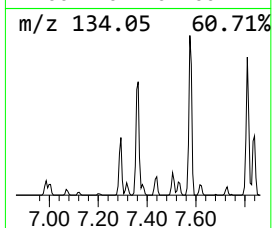
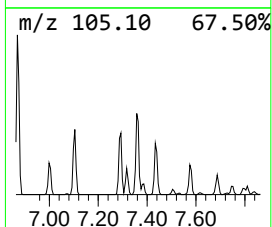
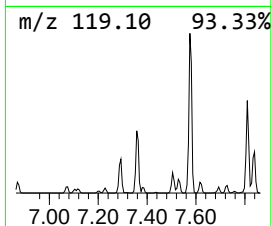
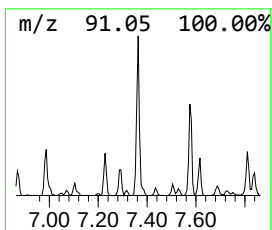
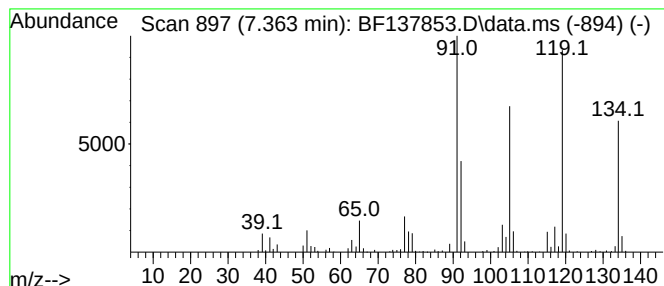
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Benzene, 1,2-diethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.363	6.10 ng	248979	1,4-Dichlorobenzene-d4	7.051

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	93
2		Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	93
3		Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93
4		1,3,8-p-Menthatriene	134	C10H14	018368-95-1	93
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

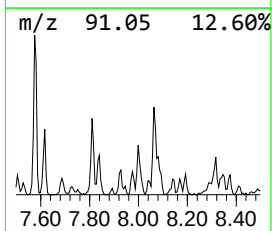
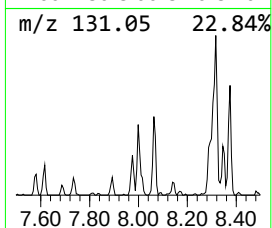
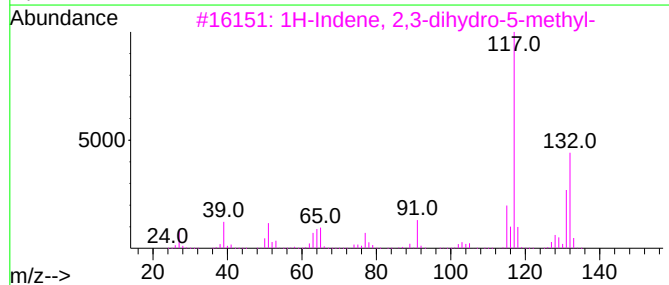
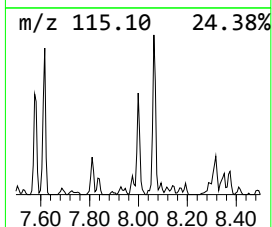
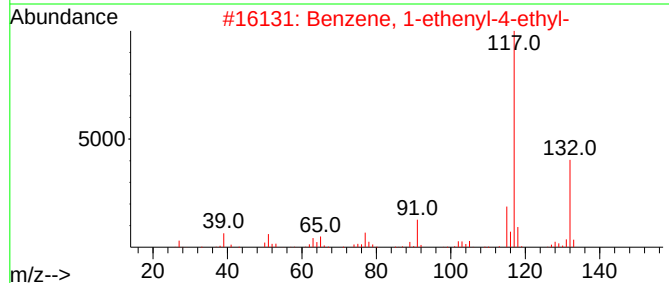
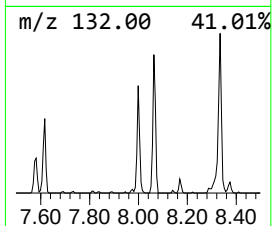
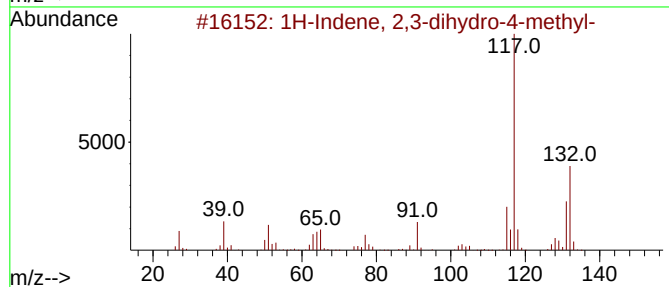
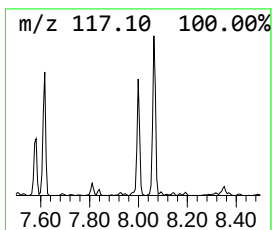
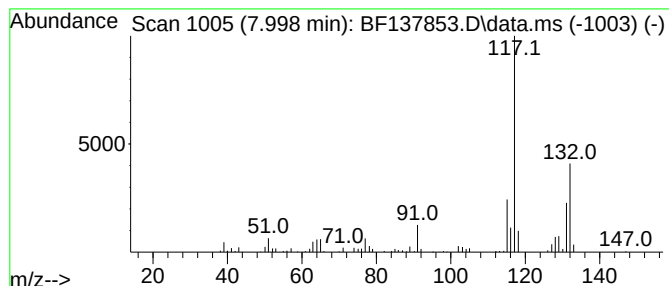
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ClientSampleId :
WC-CC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.998	2.07 ng	136805	Naphthalene-d8	8.334	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
3	1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
4	Indan, 1-methyl-	132	C10H12	000767-58-8	83
5	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
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Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

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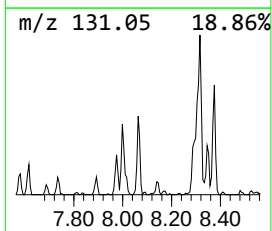
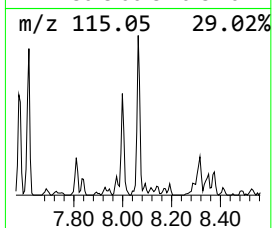
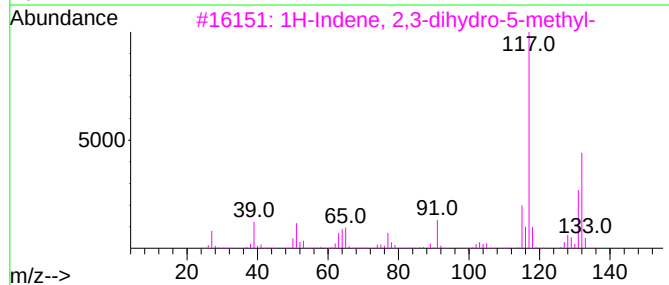
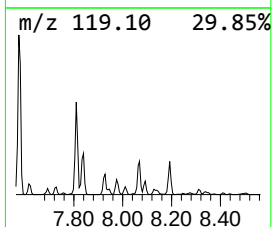
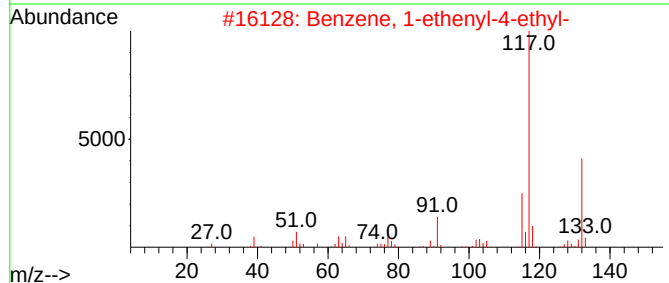
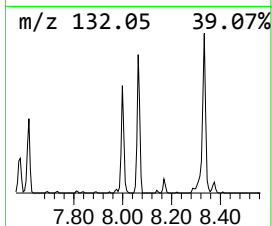
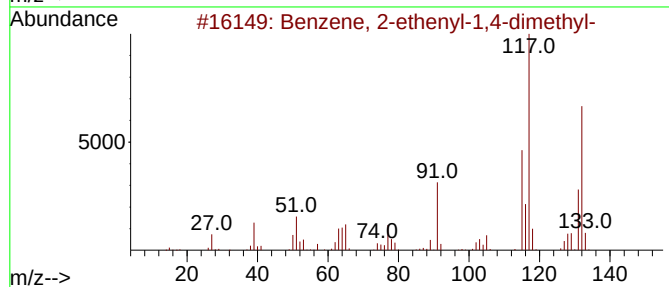
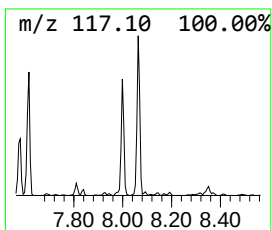
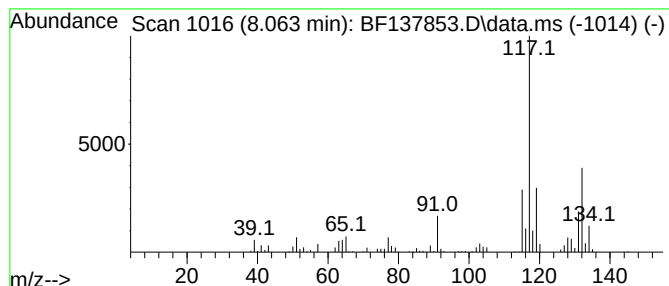
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.063	3.18 ng	209658	Naphthalene-d8	8.334

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	96
2		Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	90
3		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	90
4		3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
5		Benzene, 2-butenyl-	132	C10H12	001560-06-1	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-CC

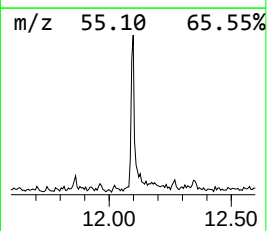
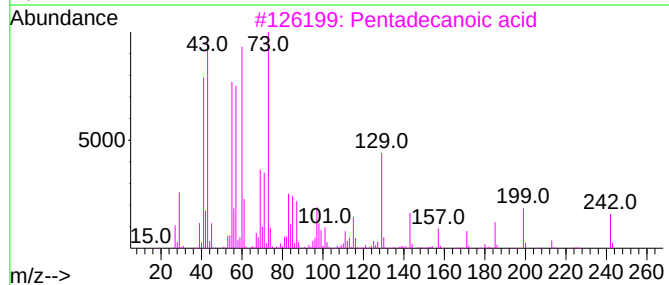
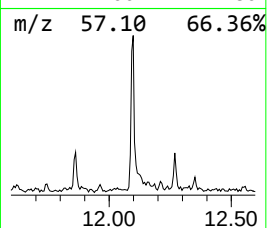
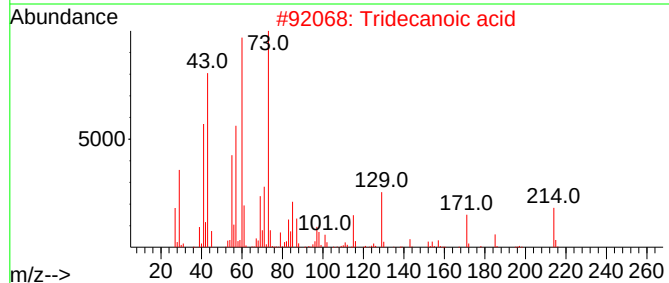
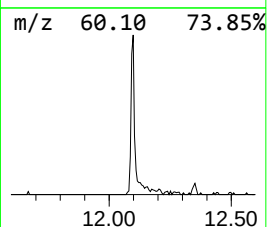
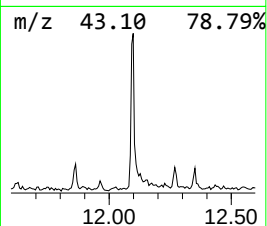
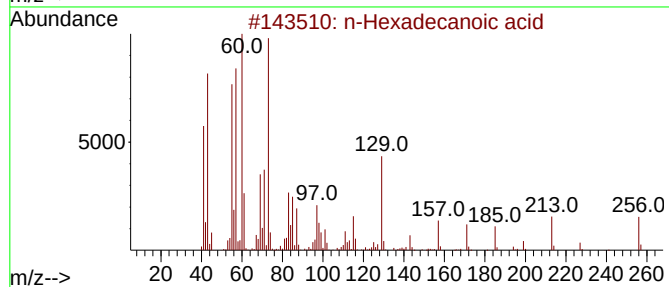
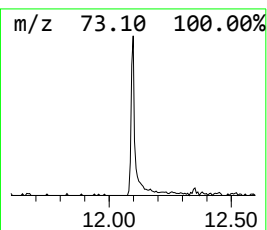
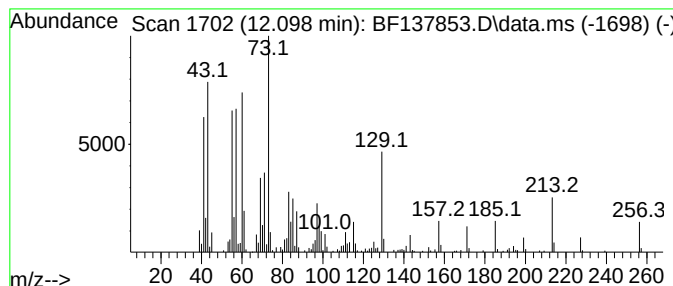
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.098	3.11 ng	206444	Phenanthrene-d10	11.592

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	91
3		Pentadecanoic acid	242	C15H30O2	001002-84-2	81
4		Tetradecanoic acid	228	C14H28O2	000544-63-8	72
5		n-Decanoic acid	172	C10H20O2	000334-48-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-CC

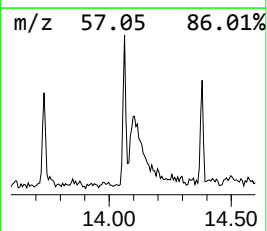
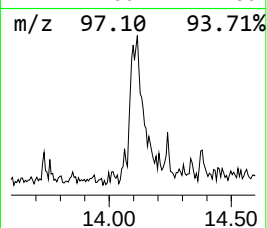
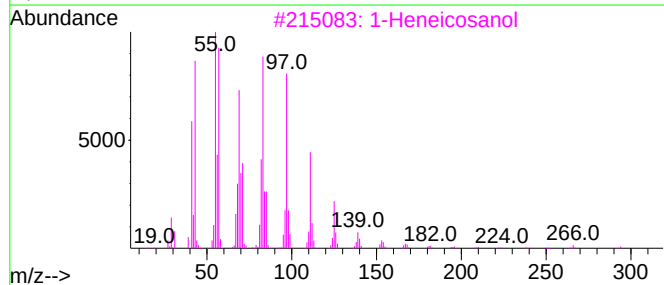
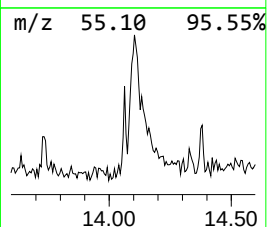
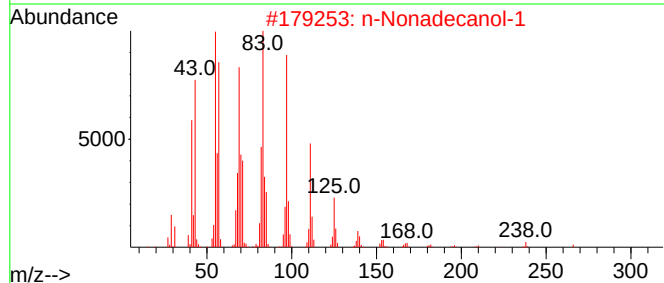
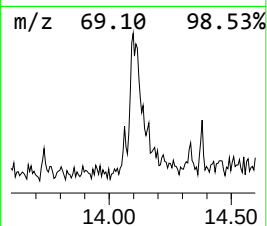
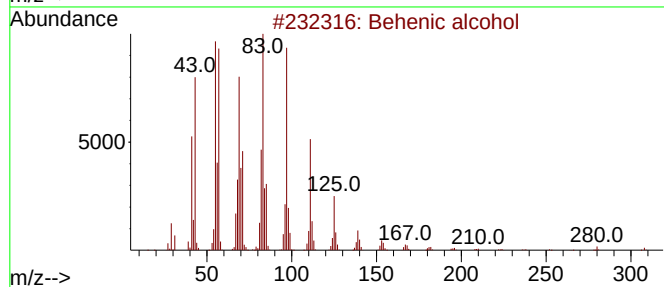
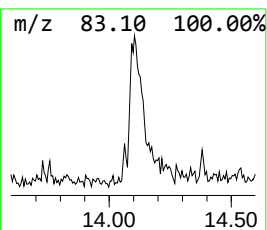
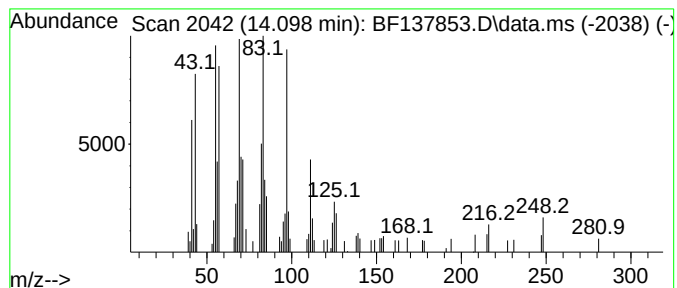
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Behenic alcohol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.098	3.91 ng	165520	Chrysene-d12	14.239

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Behenic alcohol	326	C22H46O	000661-19-8	95
2			n-Nonadecanol-1	284	C19H40O	001454-84-8	95
3			1-Heneicosanol	312	C21H44O	015594-90-8	95
4			n-Heptadecanol-1	256	C17H36O	001454-85-9	95
5			Trifluoroacetic acid, pentadecyl...	324	C17H31F3O2	959010-23-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF050824\
Data File : BF137853.D
Acq On : 08 May 2024 14:12
Operator : CG\JU
Sample : P2413-01
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
WC-CC

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF043024.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...	2.428	58.1	ng	2370720	1	7.051	816251	20.0
Heptane, 3-methyl-	4.322	2.9	ng	119600	1	7.051	816251	20.0
Heptane, 2,5-di...	5.222	2.3	ng	94195	1	7.051	816251	20.0
Octane, 4-methyl-	5.563	2.6	ng	106456	1	7.051	816251	20.0
Benzene, propyl-	6.510	5.8	ng	237998	1	7.051	816251	20.0
Mesitylene	6.869	3.6	ng	147362	1	7.051	816251	20.0
Indane	7.228	3.8	ng	153458	1	7.051	816251	20.0
Benzene, 1,4-di...	7.293	2.4	ng	98596	1	7.051	816251	20.0
Benzene, 1,2-di...	7.363	6.1	ng	248979	1	7.051	816251	20.0
1H-Indene, 2,3-...	7.998	2.1	ng	136805	2	8.334	1320500	20.0
Benzene, 2-ethe...	8.063	3.2	ng	209658	2	8.334	1320500	20.0
n-Hexadecanoic ...	12.098	3.1	ng	206444	4	11.592	1328820	20.0
Behenic alcohol	14.098	3.9	ng	165520	5	14.239	846174	20.0