

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
 Data File : BP007315.D
 Acq On : 04 Oct 2021 20:54
 Operator : CG/JU
 Sample : M4037-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 HD-01-100121

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_P\Methods\8270E-BP092121.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BP007315.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.446	394	401	411	rBV	1672923	2505360	29.32%	4.172%
2	7.046	666	673	689	rBV	1497235	2464868	28.85%	4.104%
3	7.863	805	812	822	rBV	1102255	1718470	20.11%	2.861%
4	9.028	1003	1010	1025	rBV	910888	1576737	18.45%	2.625%
5	10.451	1245	1252	1269	rBV	211772	476367	5.58%	0.793%
6	10.663	1277	1288	1303	rBV	1281560	2236982	26.18%	3.725%
7	13.122	1700	1706	1717	rBV	2291585	3436044	40.21%	5.721%
8	13.322	1736	1740	1748	rVB	69446	105466	1.23%	0.176%
9	13.981	1845	1852	1855	rBV3	44161	87959	1.03%	0.146%
10	14.392	1918	1922	1929	rVB	161657	234105	2.74%	0.390%
11	14.498	1932	1940	1949	rBV	1671289	2536183	29.68%	4.223%
12	15.345	2080	2084	2089	rVB	225953	280069	3.28%	0.466%
13	15.757	2150	2154	2159	rVB	74520	110623	1.29%	0.184%
14	15.992	2187	2194	2202	rBV	1341059	2082019	24.37%	3.467%
15	16.210	2227	2231	2240	rBV	280879	510083	5.97%	0.849%
16	17.010	2362	2367	2371	rBV	269304	335301	3.92%	0.558%
17	17.063	2372	2376	2380	rVB2	92954	129660	1.52%	0.216%
18	17.251	2403	2408	2413	rBV	1796218	2606654	30.51%	4.340%
19	17.292	2413	2415	2421	rVB	133341	189254	2.21%	0.315%
20	17.386	2428	2431	2436	rVB	62435	88013	1.03%	0.147%
21	17.751	2489	2493	2497	rBV	278519	352196	4.12%	0.586%
22	17.792	2497	2500	2504	rVV	114631	144378	1.69%	0.240%
23	17.922	2517	2522	2527	rBV	3692513	4415133	51.67%	7.352%
24	18.139	2556	2559	2562	rBV	137497	167076	1.96%	0.278%
25	18.422	2603	2607	2610	rBV	278469	317845	3.72%	0.529%
26	19.022	2704	2709	2712	rBV	6477174	8544398	100.00%	14.228%
27	19.057	2712	2715	2719	rVB	6853441	7925772	92.76%	13.197%
28	19.180	2732	2736	2741	rBV	1944099	2204647	25.80%	3.671%
29	19.263	2746	2750	2755	rBV	466387	618023	7.23%	1.029%
30	19.539	2793	2797	2801	rBV	356439	427953	5.01%	0.713%
31	19.592	2803	2806	2808	rVV2	358420	423497	4.96%	0.705%
32	19.616	2808	2810	2818	rVB	383148	536588	6.28%	0.893%
33	19.810	2839	2843	2848	rBV	3048916	3779602	44.23%	6.294%
34	20.110	2891	2894	2897	rVB2	299468	277760	3.25%	0.463%
35	20.245	2915	2917	2921	rBV2	223663	249394	2.92%	0.415%
36	20.592	2974	2976	2980	rVB	349655	322572	3.78%	0.537%

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ClientSampleId :
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Integration Parameters: rteint.p
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % of largest Peak
Start Thrs: 0.2 Max Peaks: 100
Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
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37	21.045	3051	3053	3057	rVB	468530	468036	5.48%	0.779%
38	21.339	3099	3103	3108	rBV	2022711	2609188	30.54%	4.345%
39	23.751	3509	3513	3519	rVB2	1361715	2560923	29.97%	4.264%

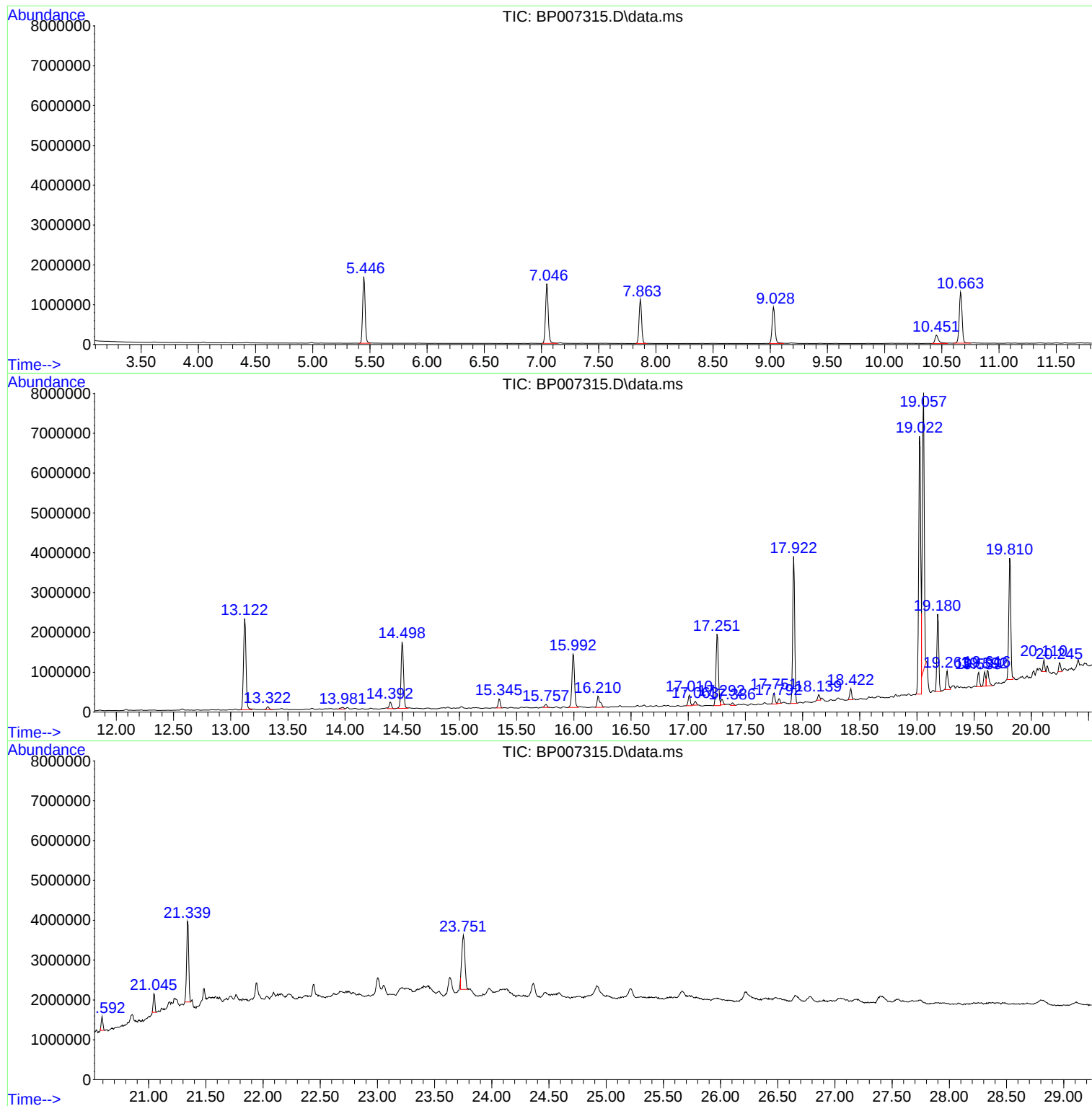
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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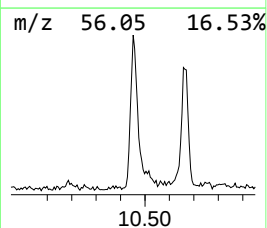
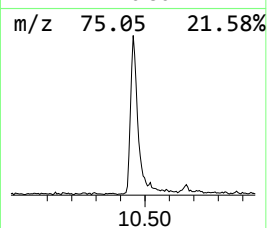
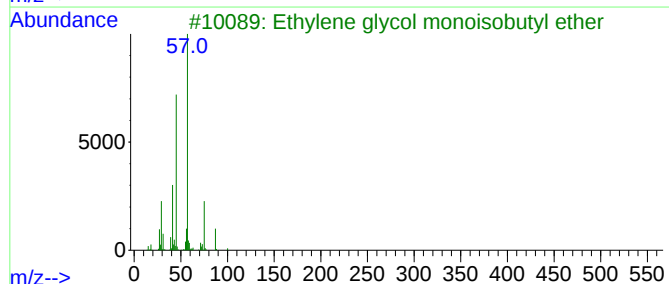
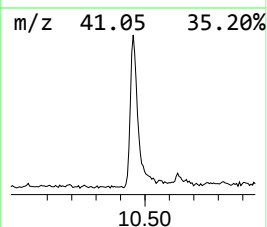
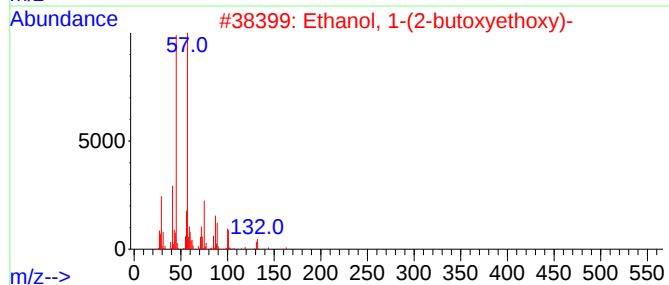
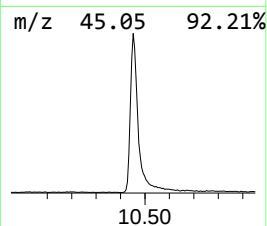
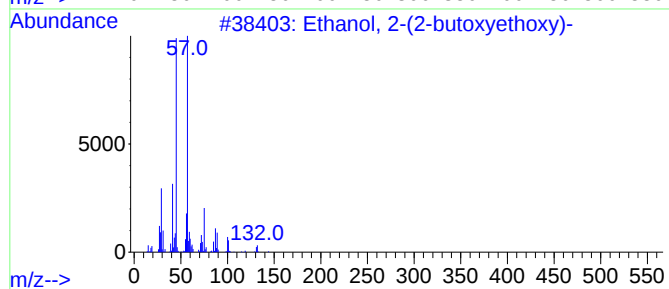
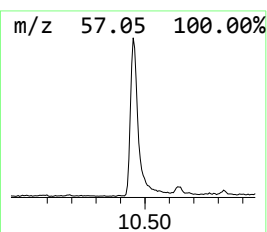
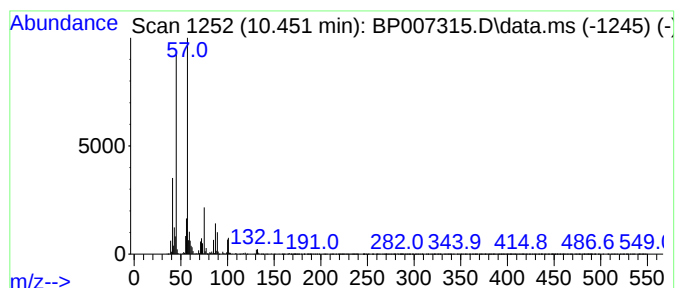
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 1 Ethanol, 2-(2-butoxyethoxy)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.451	4.26 ng	476367	Naphthalene-d8	10.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	90
2			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	83
3			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	64
4			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56
5			Tetraethylene glycol	194	C8H18O5	000112-60-7	50



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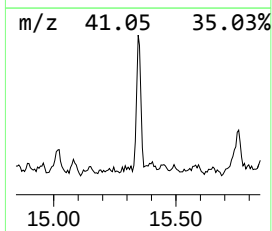
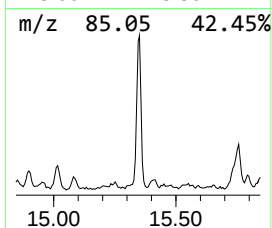
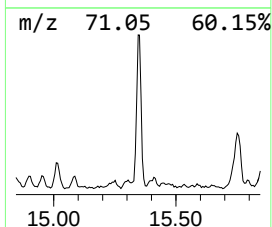
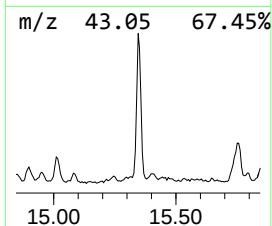
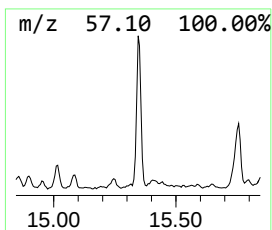
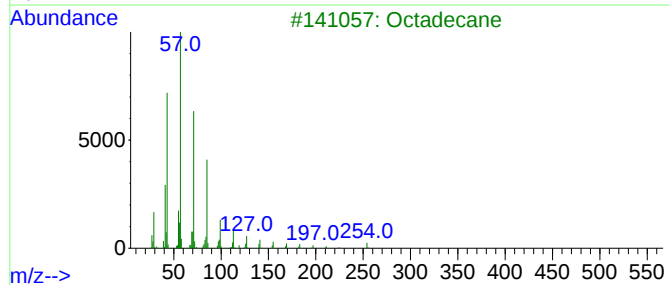
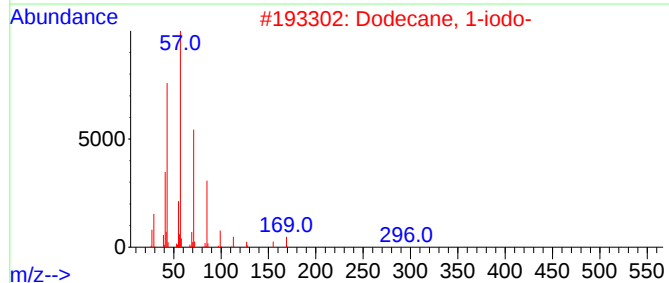
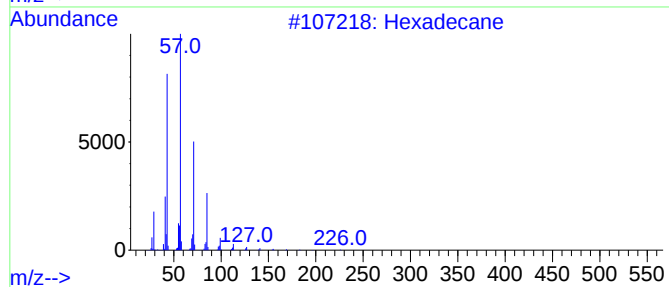
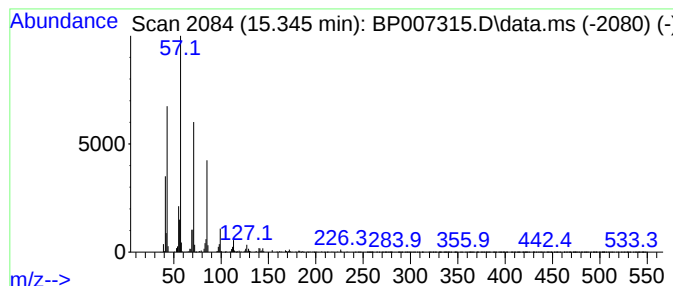
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
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TIC Library : C:\Database\NIST20.L
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Peak Number 2 Hexadecane Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.345	2.21 ng	280069	Acenaphthene-d10	14.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	97
2			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	90
3			Octadecane	254	C18H38	000593-45-3	90
4			Nonadecane	268	C19H40	000629-92-5	90
5			Tetradecane	198	C14H30	000629-59-4	90



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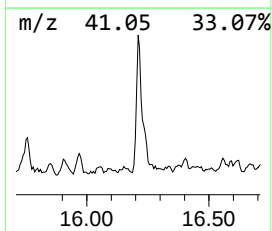
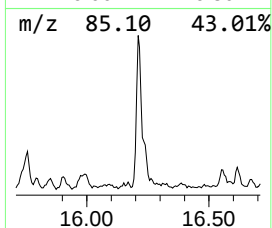
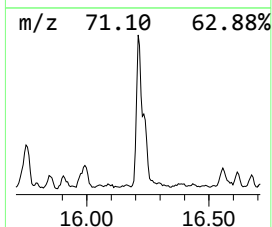
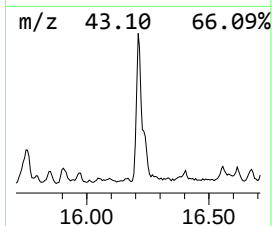
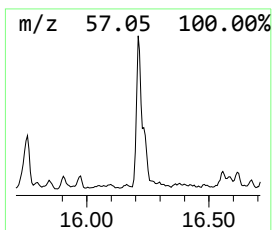
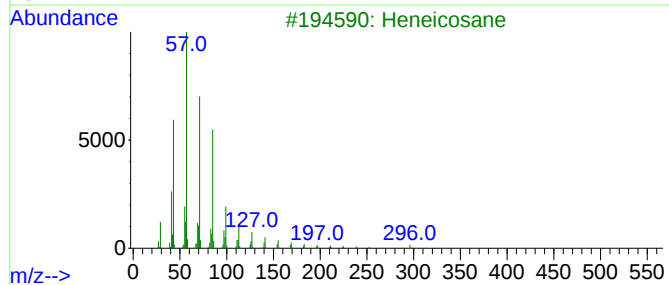
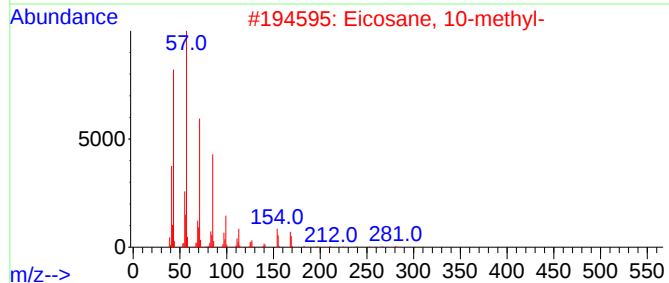
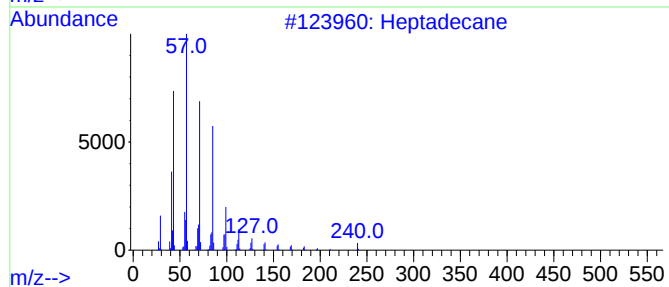
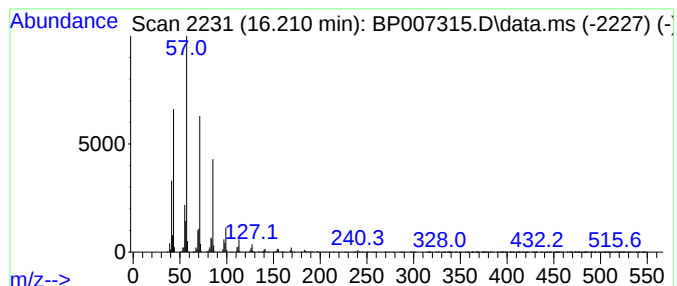
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 3 Heptadecane Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.210	3.91 ng	510083	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	96
2			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
3			Heneicosane	296	C21H44	000629-94-7	90
4			Nonadecane	268	C19H40	000629-92-5	90
5			Pentadecane	212	C15H32	000629-62-9	90



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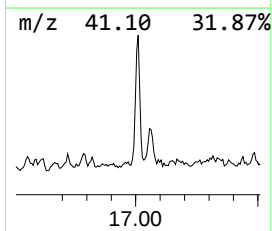
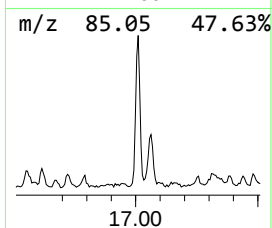
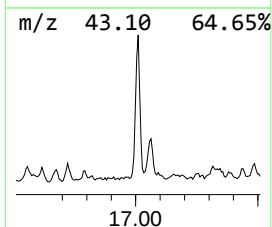
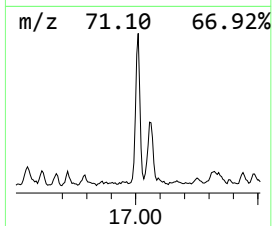
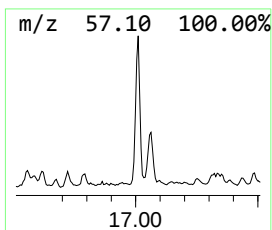
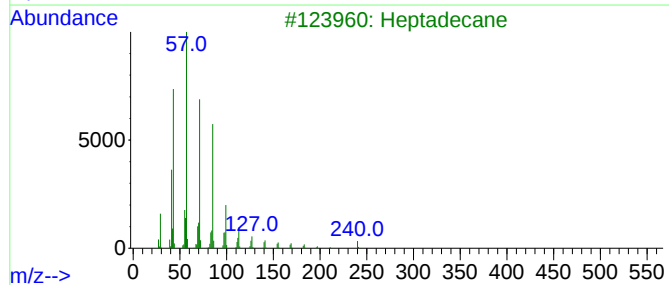
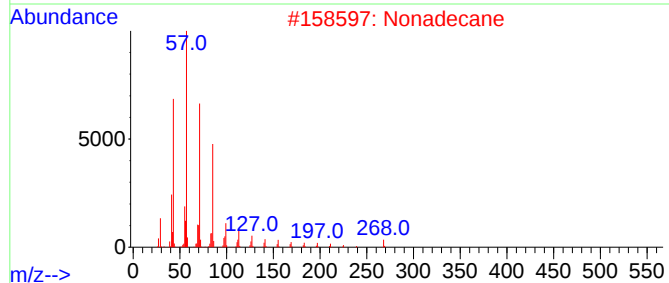
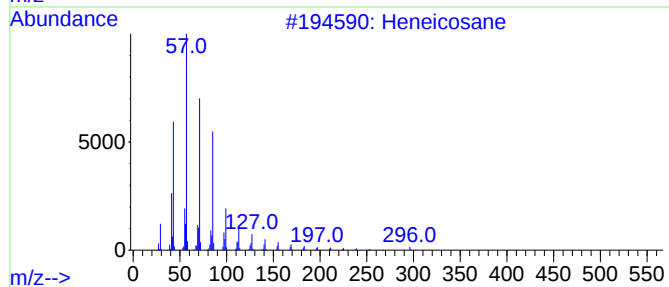
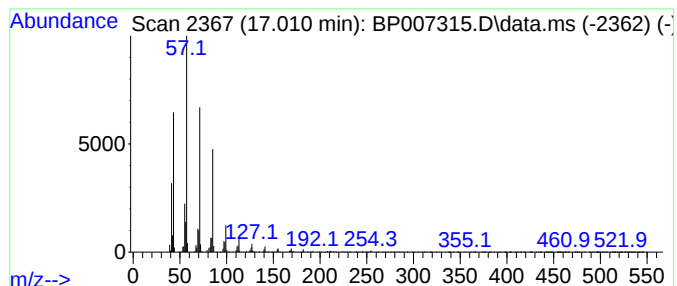
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 4 Heneicosane Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.010	2.57 ng	335301	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heneicosane	296	C21H44	000629-94-7	91
2			Nonadecane	268	C19H40	000629-92-5	91
3			Heptadecane	240	C17H36	000629-78-7	91
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	86
5			Eicosane	282	C20H42	000112-95-8	86



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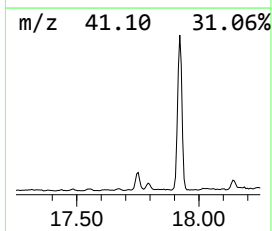
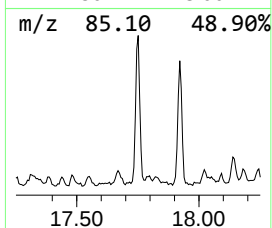
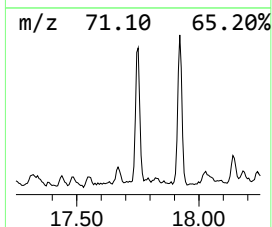
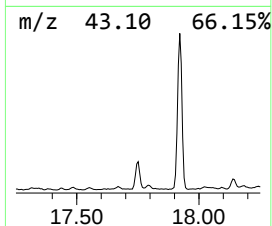
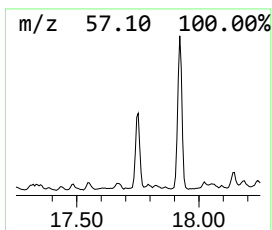
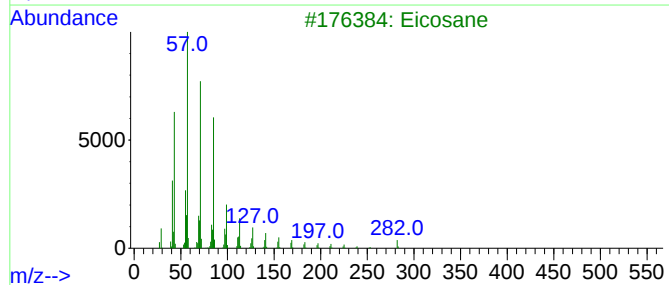
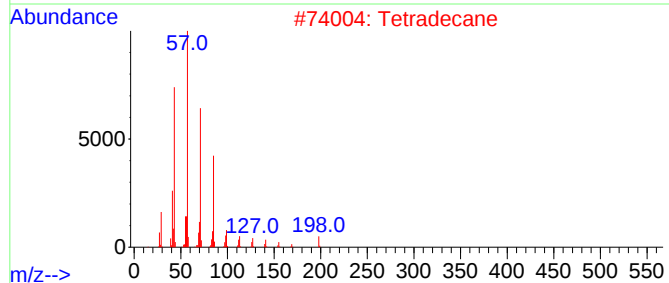
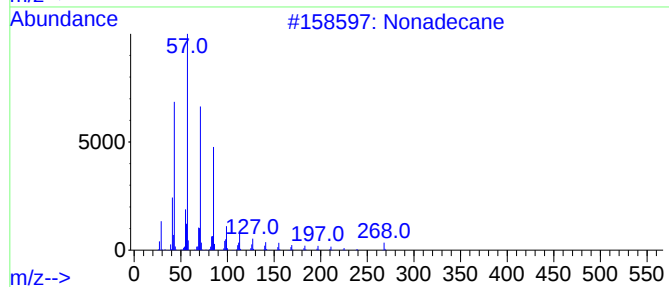
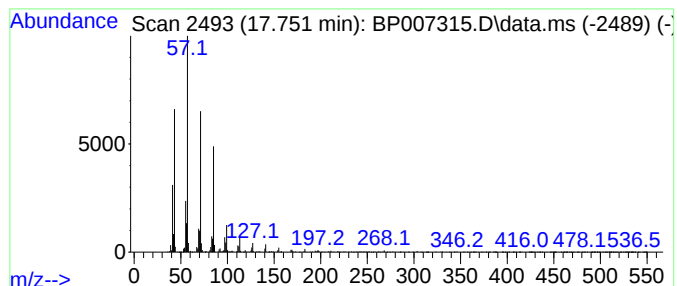
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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Peak Number 5 Nonadecane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.751	2.70 ng	352196	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	94
2			Tetradecane	198	C14H30	000629-59-4	91
3			Eicosane	282	C20H42	000112-95-8	90
4			Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	86
5			Octacosane	394	C28H58	000630-02-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

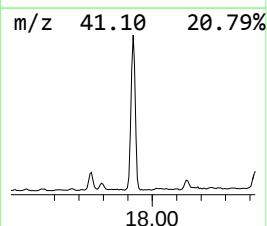
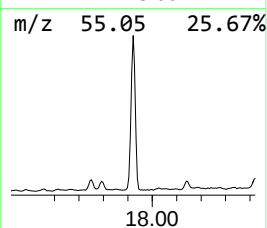
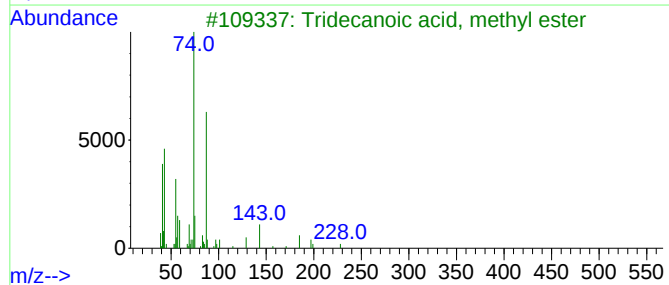
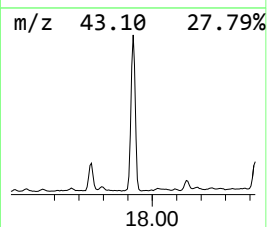
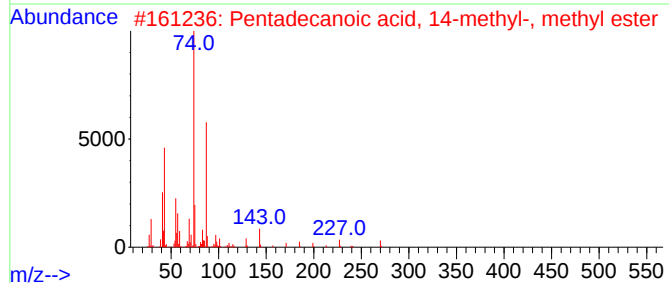
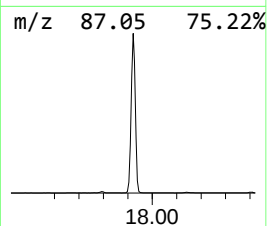
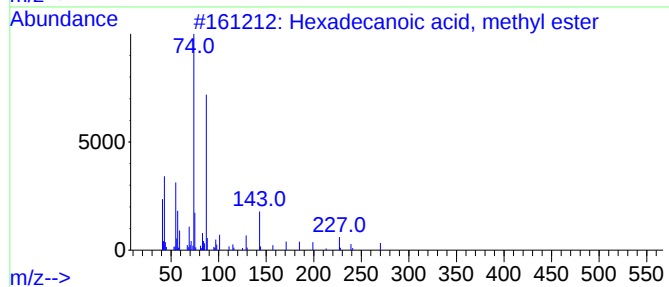
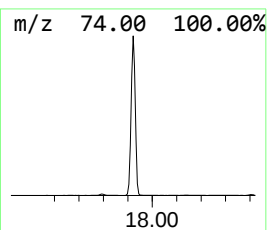
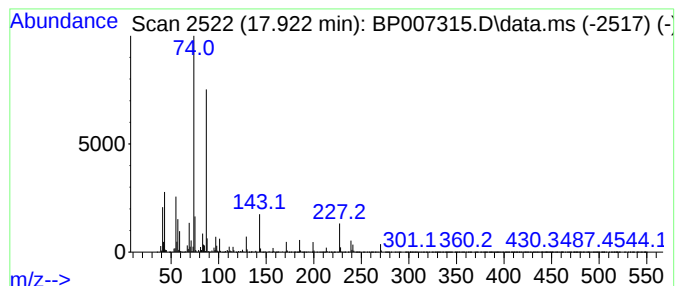
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Hexadecanoic acid, methyl e... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.922	33.88 ng	4415130	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	98
2			Pentadecanoic acid, 14-methyl-, ...	270	C17H34O2	005129-60-2	91
3			Tridecanoic acid, methyl ester	228	C14H28O2	001731-88-0	90
4			Decanoic acid, methyl ester	186	C11H22O2	000110-42-9	87
5			Methyl tetradecanoate	242	C15H30O2	000124-10-7	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

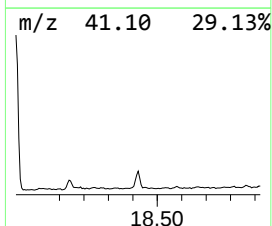
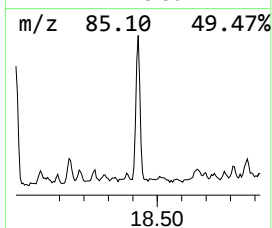
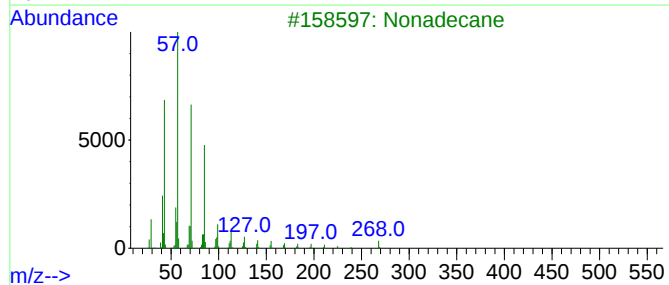
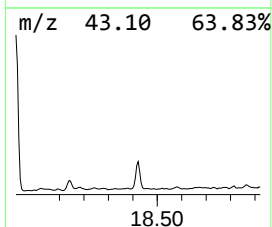
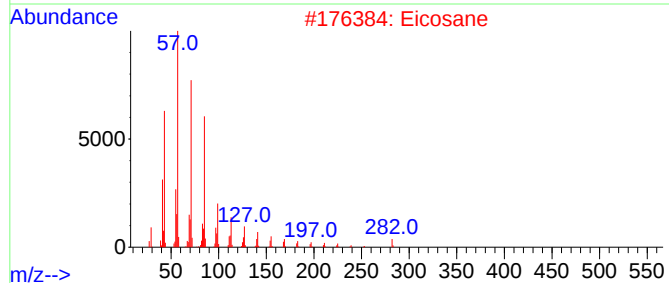
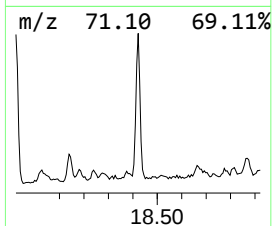
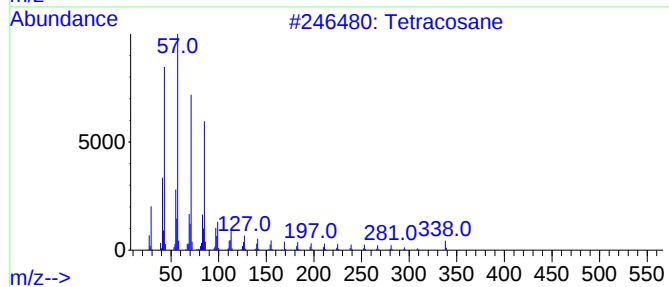
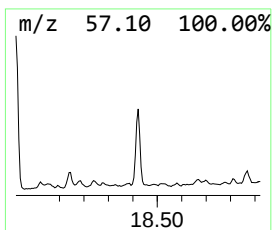
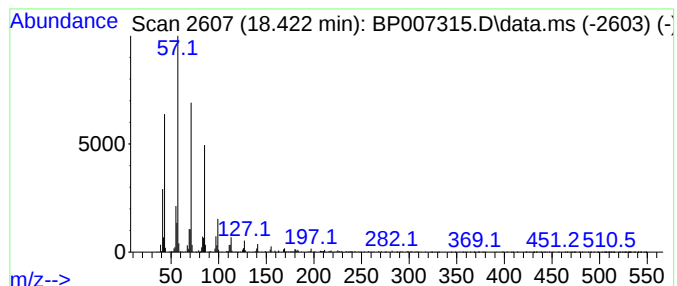
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tetracosane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.422	2.44 ng	317845	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Eicosane	282	C20H42	000112-95-8	97
3			Nonadecane	268	C19H40	000629-92-5	94
4			Dodecane, 2-methyl-6-propyl-	226	C16H34	055045-08-4	93
5			Heptadecane	240	C17H36	000629-78-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

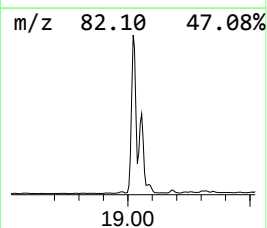
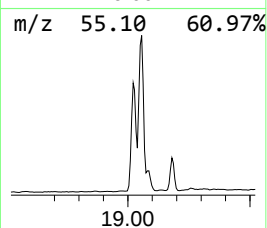
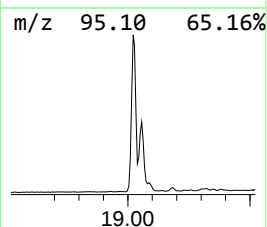
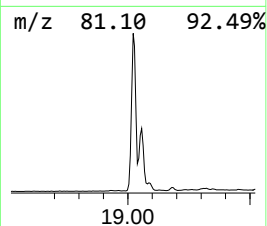
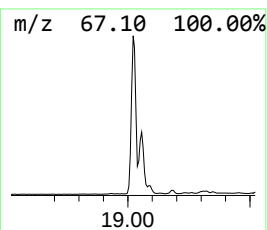
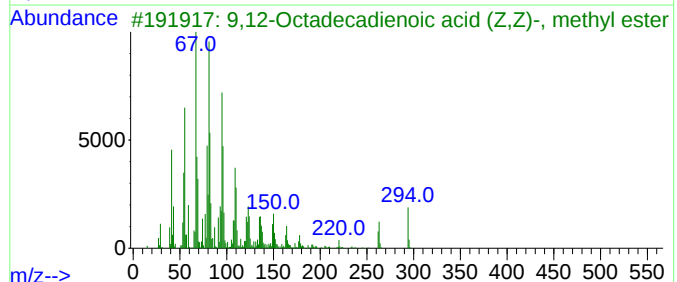
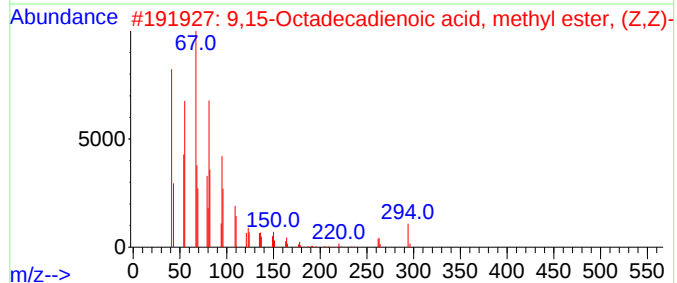
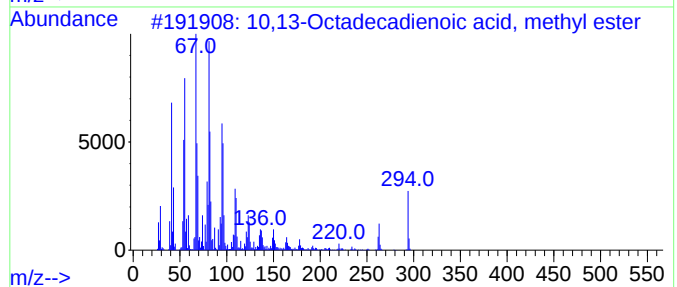
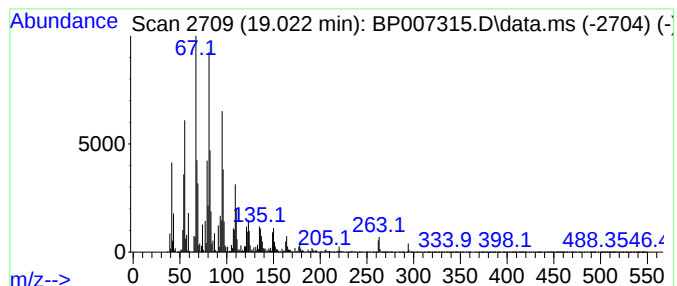
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 10,13-Octadecadienoic acid,... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.022	65.56 ng	8544400	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	99		
2	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99		
3	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99		
4	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99		
5	Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
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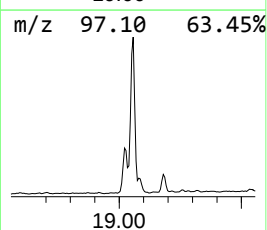
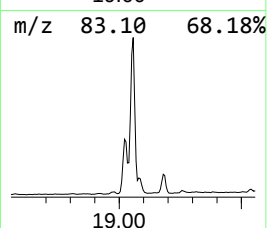
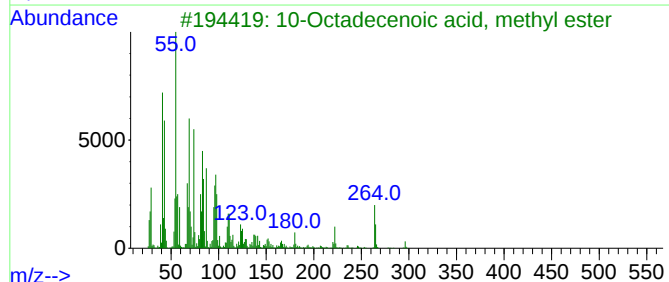
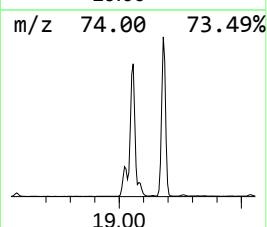
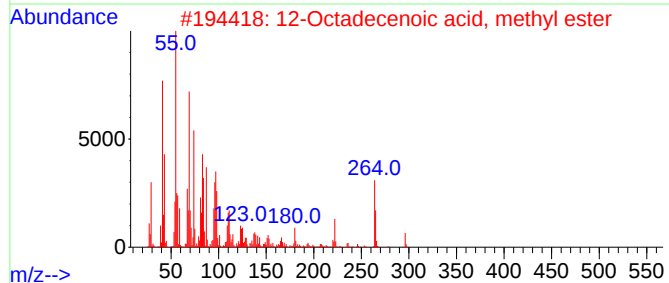
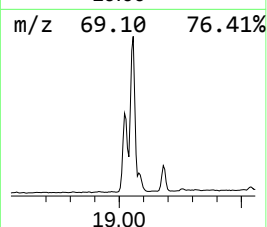
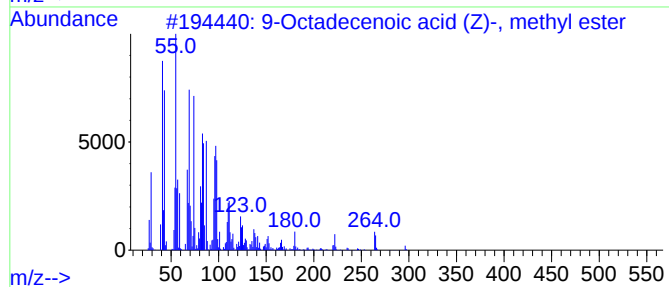
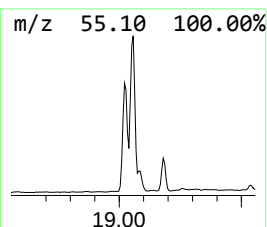
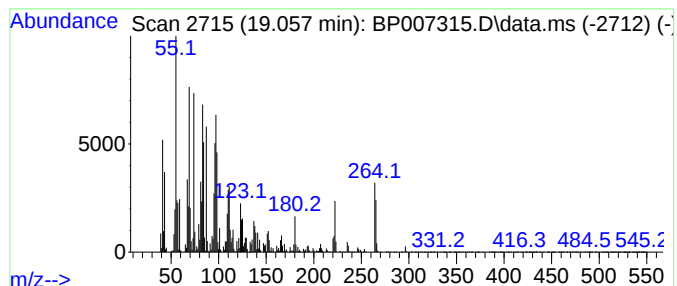
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 9-Octadecenoic acid (Z)-, m... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.057	60.81 ng	7925770	Phenanthrene-d10	17.251	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	99
2	12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	99
3	10-Octadecenoic acid, methyl ester	296	C19H36O2	013481-95-3	99
4	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	99
5	11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

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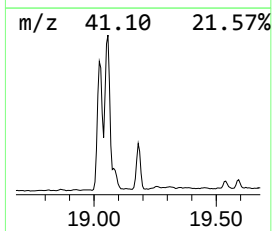
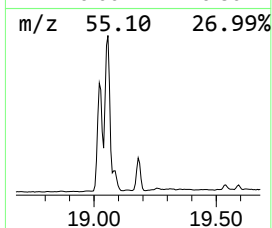
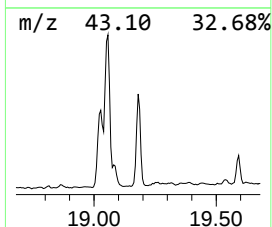
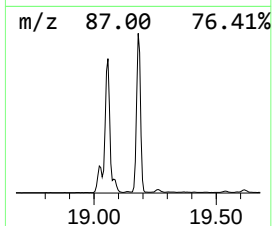
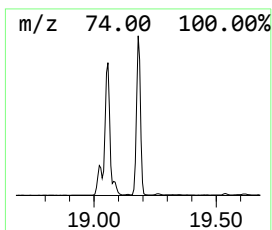
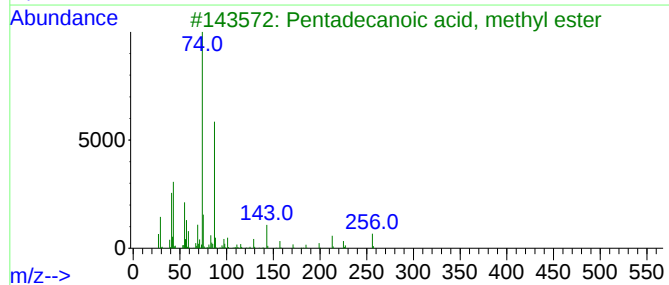
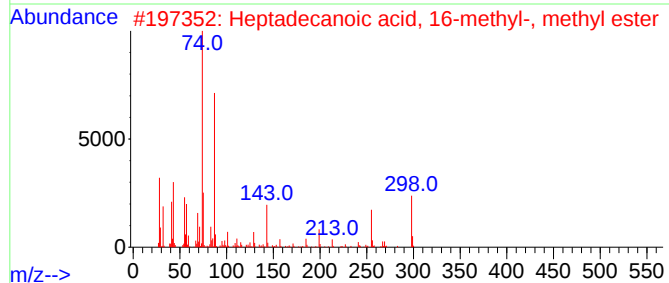
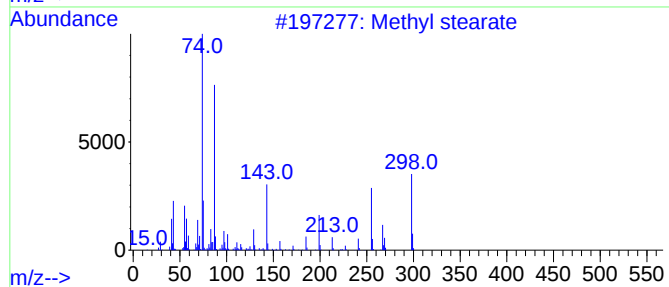
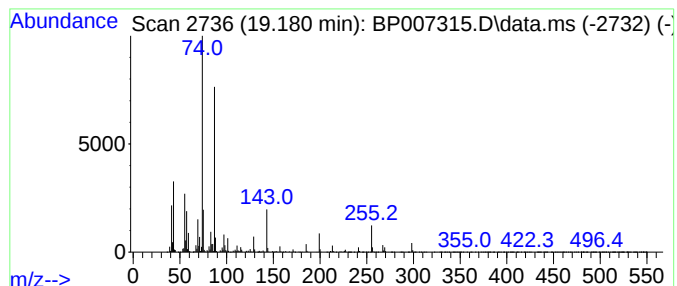
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Methyl stearate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.180	16.92 ng	2204650	Phenanthrene-d10	17.251

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl stearate	298	C19H38O2	000112-61-8	98
2			Heptadecanoic acid, 16-methyl-, ...	298	C19H38O2	005129-61-3	90
3			Pentadecanoic acid, methyl ester	256	C16H32O2	007132-64-1	87
4			Hexadecanoic acid, methyl ester	270	C17H34O2	000112-39-0	86
5			Heptadecanoic acid, 15-methyl-, ...	298	C19H38O2	054833-55-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
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Misc :
ALS Vial : 17 Sample Multiplier: 1

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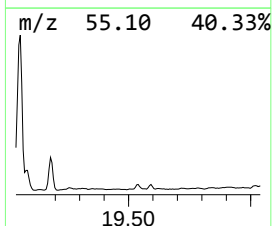
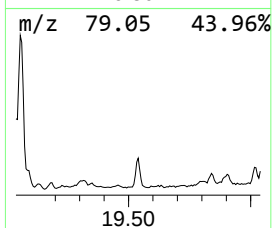
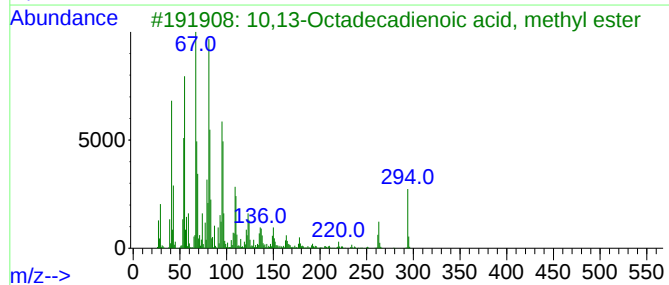
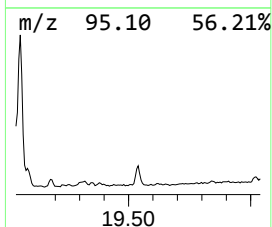
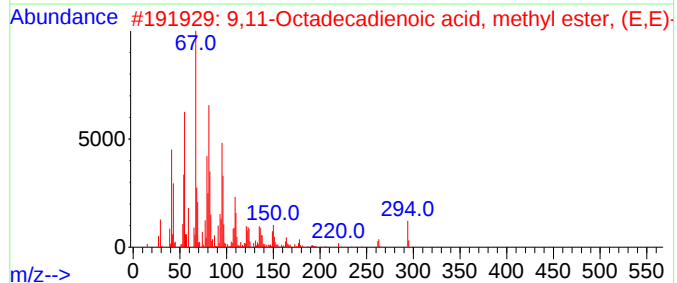
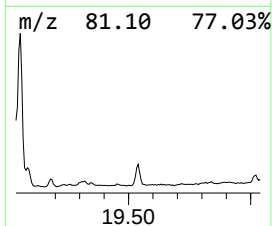
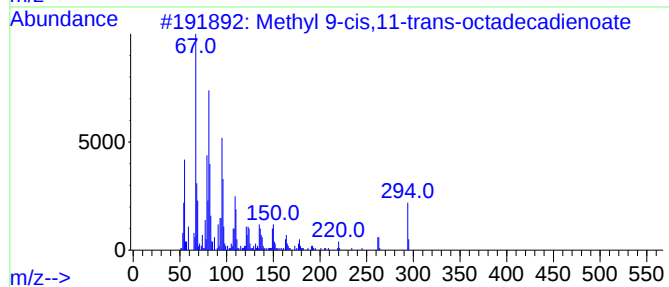
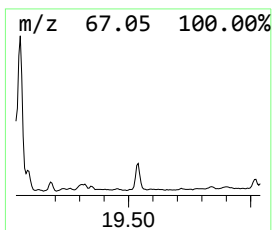
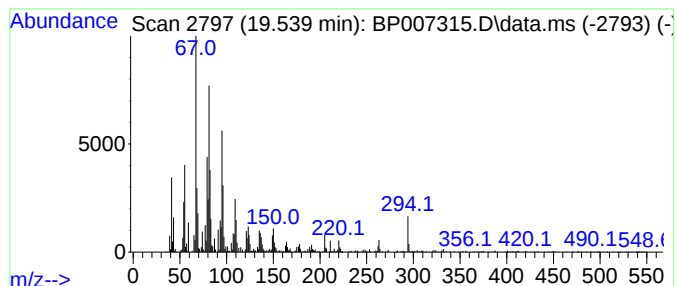
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Methyl 9-cis,11-trans-octad... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	3.28 ng	427953	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 9-cis,11-trans-octadecadi...	294	C19H34O2	013058-52-1	97
2			9,11-Octadecadienoic acid, methy...	294	C19H34O2	013038-47-6	96
3			10,13-Octadecadienoic acid, meth...	294	C19H34O2	056554-62-2	93
4			6-Tetradecyne	194	C14H26	003730-08-3	93
5			Methyl 10-trans,12-cis-octadecad...	294	C19H34O2	1000336-44-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
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Operator : CG/JU
Sample : M4037-01 2X
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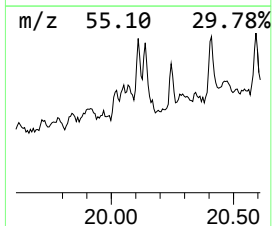
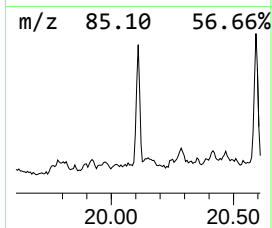
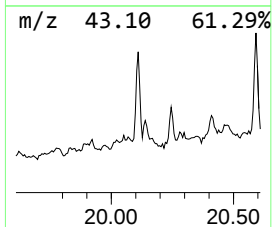
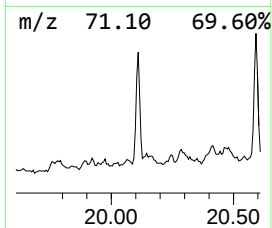
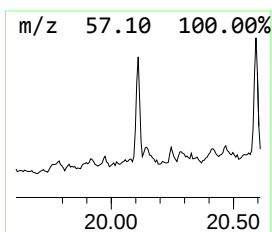
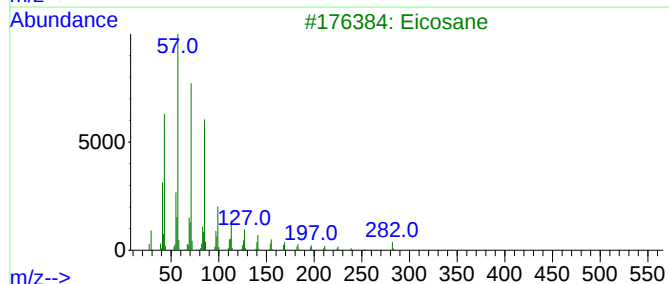
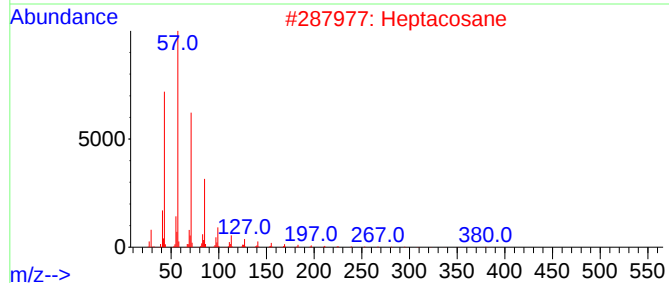
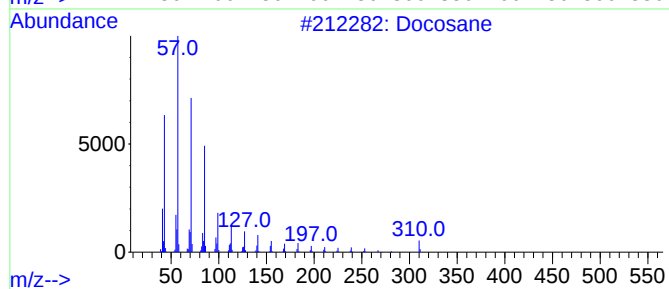
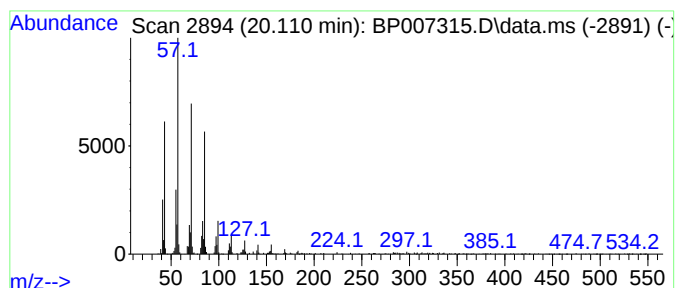
Instrument :
BNA_P
ClientSampleId :
HD-01-100121

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Docosane Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.110	2.13 ng	277760	Chrysene-d12	21.339	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane	310	C22H46	000629-97-0	95
2	Heptacosane	380	C27H56	000593-49-7	93
3	Eicosane	282	C20H42	000112-95-8	90
4	Heneicosane	296	C21H44	000629-94-7	87
5	Hexacosane	366	C26H54	000630-01-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

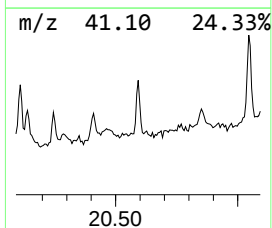
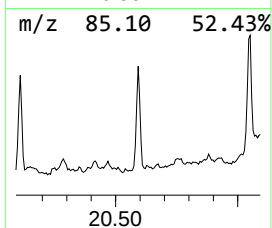
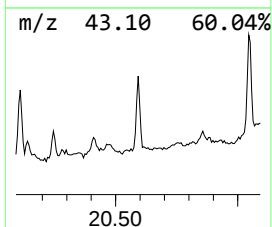
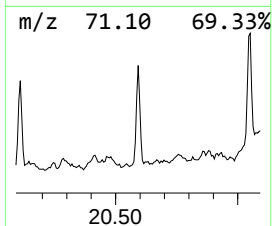
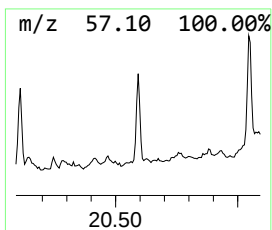
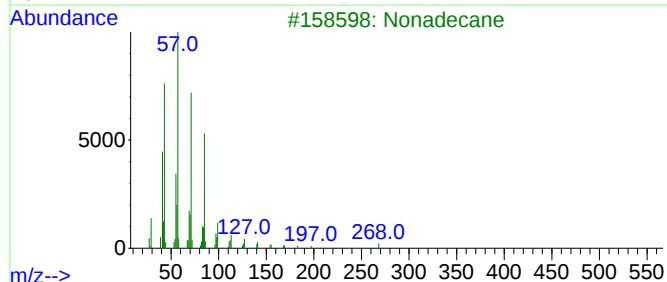
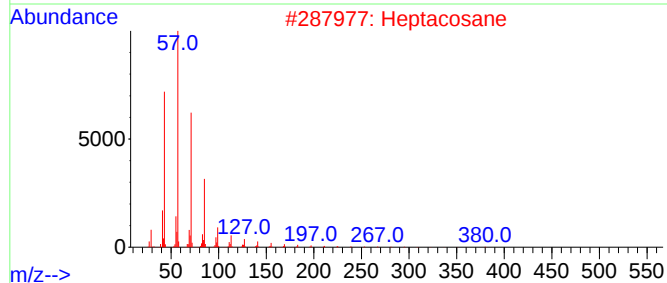
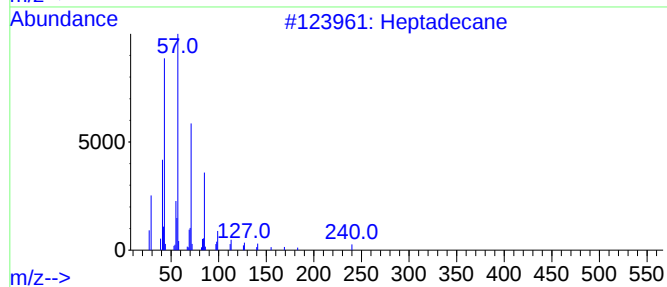
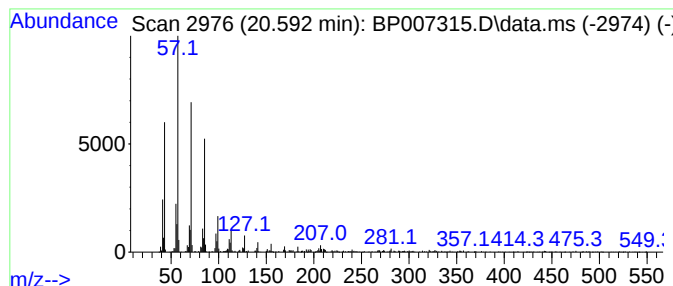
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptacosane Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.592	2.47 ng	322572	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	95
2			Heptacosane	380	C27H56	000593-49-7	94
3			Nonadecane	268	C19H40	000629-92-5	93
4			Hexacosane	366	C26H54	000630-01-3	90
5			Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

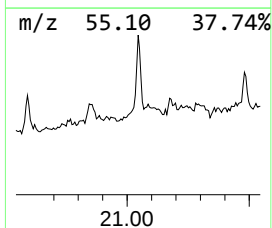
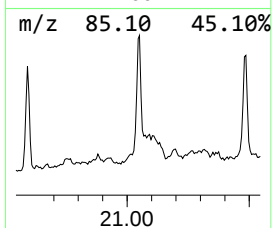
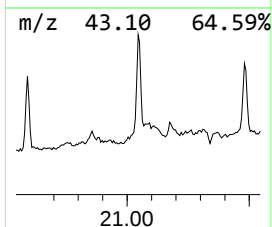
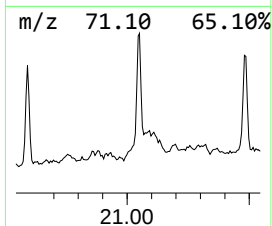
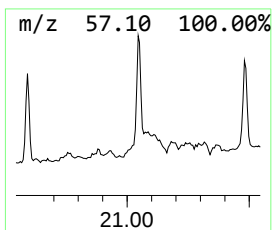
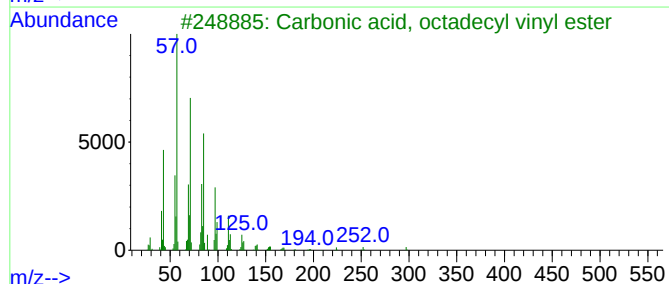
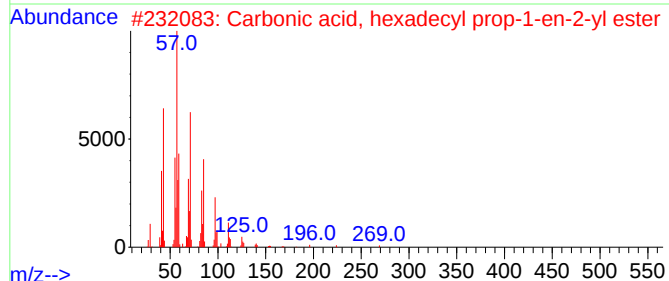
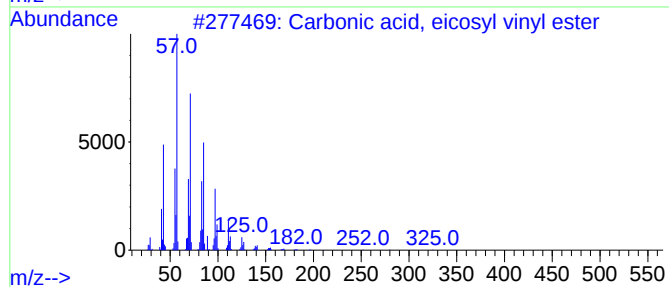
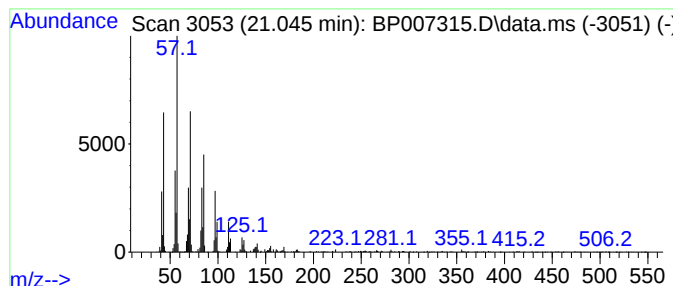
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 14 Carbonic acid, eicosyl vinyl... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.045	3.59 ng	468036	Chrysene-d12	21.339

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	94
2			Carbonic acid, hexadecyl prop-1-...	326	C20H38O3	1000382-90-3	91
3			Carbonic acid, octadecyl vinyl e...	340	C21H40O3	1000382-54-4	91
4			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	91
5			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100421\
Data File : BP007315.D
Acq On : 04 Oct 2021 20:54
Operator : CG/JU
Sample : M4037-01 2X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
HD-01-100121

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP092121.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
Ethanol, 2-(2-b...	10.451	4.3	ng	476367	2	10.663	2236980	20.0
Hexadecane	15.345	2.2	ng	280069	3	14.498	2536180	20.0
Heptadecane	16.210	3.9	ng	510083	4	17.251	2606650	20.0
Heneicosane	17.010	2.6	ng	335301	4	17.251	2606650	20.0
Nonadecane	17.751	2.7	ng	352196	4	17.251	2606650	20.0
Hexadecanoic ac...	17.922	33.9	ng	4415130	4	17.251	2606650	20.0
Tetracosane	18.422	2.4	ng	317845	4	17.251	2606650	20.0
10,13-Octadecad...	19.022	65.6	ng	8544400	4	17.251	2606650	20.0
9-Octadecenoic ...	19.057	60.8	ng	7925770	4	17.251	2606650	20.0
Methyl stearate	19.180	16.9	ng	2204650	4	17.251	2606650	20.0
Methyl 9-cis,11...	19.539	3.3	ng	427953	5	21.339	2609190	20.0
Docosane	20.110	2.1	ng	277760	5	21.339	2609190	20.0
Heptacosane	20.592	2.5	ng	322572	5	21.339	2609190	20.0
Carbonic acid, ...	21.045	3.6	ng	468036	5	21.339	2609190	20.0