

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
 Data File : BF125955.D
 Acq On : 27 Oct 2021 2:30
 Operator : JU/CG
 Sample : M4355-01 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-2-102521

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-BF102521.M

Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF125955.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.128	3	7	12	rVB	1411775	1711484	83.56%	7.650%
2	3.934	310	314	317	rVB	26775	33669	1.64%	0.150%
3	4.516	408	413	419	rBV3	12805	24282	1.19%	0.109%
4	5.075	505	508	513	rVB	19672	21908	1.07%	0.098%
5	5.475	571	576	579	rBV	2269112	2048246	100.00%	9.155%
6	6.481	743	747	750	rBV	2274538	1947947	95.10%	8.707%
7	6.869	809	813	816	rBV	1377839	1181174	57.67%	5.279%
8	7.422	903	907	910	rBV	1544351	1193532	58.27%	5.335%
9	7.510	918	922	927	rVB2	22197	23946	1.17%	0.107%
10	8.051	1010	1014	1017	rBV	380933	265923	12.98%	1.189%
11	8.151	1026	1031	1034	rBV	1670561	1433601	69.99%	6.408%
12	8.575	1100	1103	1106	rBV	25835	22150	1.08%	0.099%
13	8.745	1129	1132	1136	rBV	76516	59693	2.91%	0.267%
14	8.957	1161	1168	1173	rBV6	17794	38629	1.89%	0.173%
15	9.039	1178	1182	1185	rBV6	15564	29419	1.44%	0.131%
16	9.110	1190	1194	1197	rVB3	20929	23910	1.17%	0.107%
17	9.175	1203	1205	1210	rVB3	37583	28790	1.41%	0.129%
18	9.222	1210	1213	1217	rBV	2073245	1650536	80.58%	7.377%
19	9.310	1225	1228	1233	rBV	118261	106572	5.20%	0.476%
20	9.375	1237	1239	1244	rVB5	29143	31228	1.52%	0.140%
21	9.480	1252	1257	1262	rBV8	19169	37615	1.84%	0.168%
22	9.563	1269	1271	1273	rBV2	25933	22313	1.09%	0.100%
23	9.586	1273	1275	1278	rVB2	35425	25403	1.24%	0.114%
24	9.633	1278	1283	1285	rBV2	65627	78440	3.83%	0.351%
25	9.822	1310	1315	1316	rBV4	22896	25972	1.27%	0.116%
26	9.839	1316	1318	1321	rVV	150642	104408	5.10%	0.467%
27	9.904	1325	1329	1332	rVV	1590407	1268732	61.94%	5.671%
28	9.933	1332	1334	1339	rVB3	27399	25876	1.26%	0.116%
29	10.016	1345	1348	1351	rVB4	25109	24357	1.19%	0.109%
30	10.080	1354	1359	1360	rVV5	20404	29737	1.45%	0.133%
31	10.104	1360	1363	1365	rVB3	41794	41401	2.02%	0.185%
32	10.163	1371	1373	1376	rVB3	28401	27451	1.34%	0.123%
33	10.316	1396	1399	1400	rBV2	24493	27275	1.33%	0.122%
34	10.339	1400	1403	1406	rVB	170496	130334	6.36%	0.583%
35	10.445	1418	1421	1424	rBV3	35432	32191	1.57%	0.144%
36	10.563	1435	1441	1446	rBV	91541	129837	6.34%	0.580%

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Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

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37	10.610	1446	1449	1453	rVB6	28063	31147	1.52%	0.139%
38	10.651	1453	1456	1458	rBV3	23224	31369	1.53%	0.140%
39	10.686	1458	1462	1465	rBV	1128023	946940	46.23%	4.232%
40	10.810	1480	1483	1492	rVB	180408	245993	12.01%	1.099%
41	11.263	1557	1560	1562	rBV	164356	127025	6.20%	0.568%
42	11.292	1562	1565	1570	rVB2	125928	129553	6.33%	0.579%
43	11.386	1577	1581	1583	rBV	1524025	1236235	60.36%	5.525%
44	11.410	1583	1585	1588	rVV	241437	182120	8.89%	0.814%
45	11.439	1588	1590	1591	rVV2	33410	28956	1.41%	0.129%
46	11.457	1591	1593	1597	rVB	48409	48434	2.36%	0.216%
47	11.569	1610	1612	1615	rVB4	24807	23484	1.15%	0.105%
48	11.610	1617	1619	1622	rBV3	27182	30071	1.47%	0.134%
49	11.686	1630	1632	1635	rBV	148780	113973	5.56%	0.509%
50	11.716	1635	1637	1639	rVB2	29838	24923	1.22%	0.111%
51	11.916	1668	1671	1675	rVB	186249	170772	8.34%	0.763%
52	11.998	1681	1685	1687	rBV2	83582	98556	4.81%	0.441%
53	12.098	1699	1702	1704	rBV	146090	123200	6.01%	0.551%
54	12.121	1704	1706	1709	rVB2	39455	35484	1.73%	0.159%
55	12.316	1736	1739	1745	rBV5	45989	85275	4.16%	0.381%
56	12.439	1757	1760	1763	rBV2	78442	81437	3.98%	0.364%
57	12.486	1765	1768	1771	rVB	149346	154720	7.55%	0.692%
58	12.598	1784	1787	1790	rBV	394571	324587	15.85%	1.451%
59	12.704	1802	1805	1807	rBV3	99456	101364	4.95%	0.453%
60	12.827	1823	1826	1828	rVV	272731	208024	10.16%	0.930%
61	12.857	1829	1831	1834	rVB	158564	123607	6.03%	0.552%
62	12.974	1847	1851	1854	rBV	1653964	1444760	70.54%	6.457%
63	13.157	1880	1882	1885	rBV2	73166	63023	3.08%	0.282%
64	13.215	1889	1892	1895	rVV2	191351	173142	8.45%	0.774%
65	13.557	1947	1950	1953	rBV	218675	203043	9.91%	0.908%
66	14.027	2026	2030	2032	rBV	950526	911094	44.48%	4.072%
67	14.204	2058	2060	2063	rVB	214735	174419	8.52%	0.780%
68	14.510	2110	2112	2115	rBV	223655	189842	9.27%	0.849%
69	15.498	2277	2280	2284	rVB2	519014	598814	29.24%	2.676%

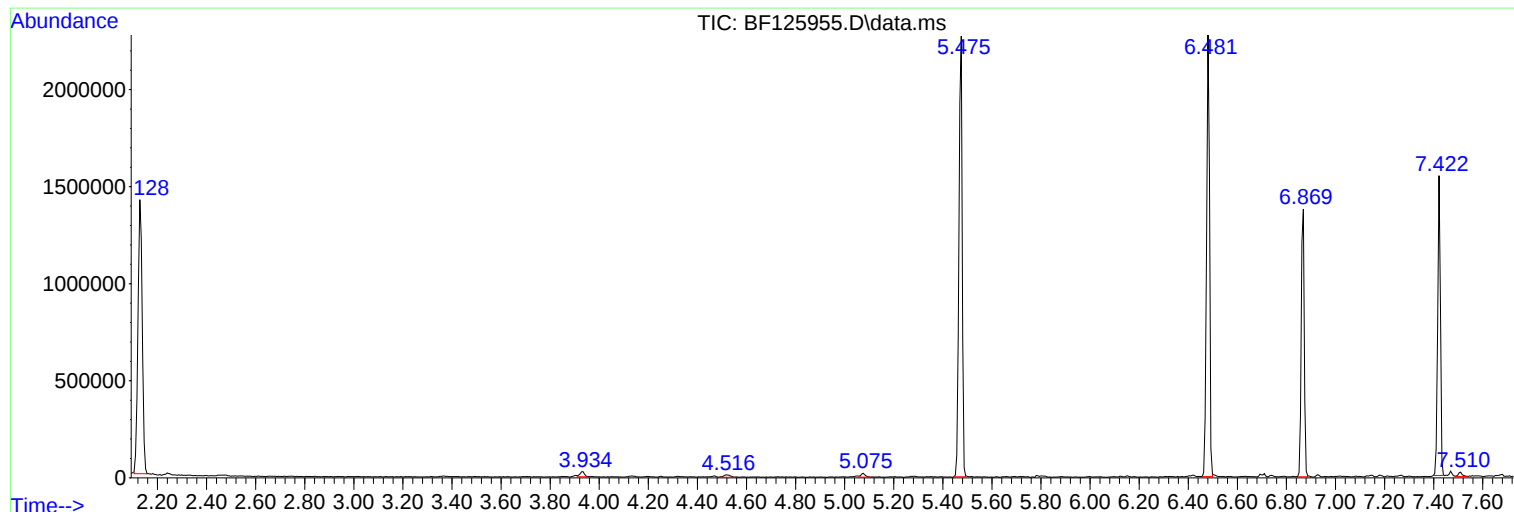
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ClientSampleId :
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Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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



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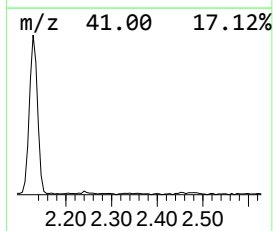
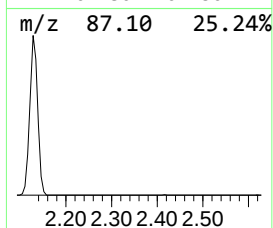
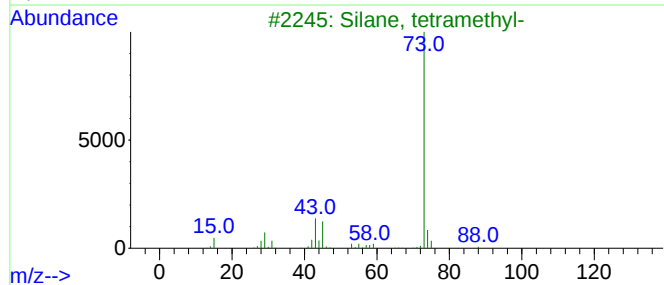
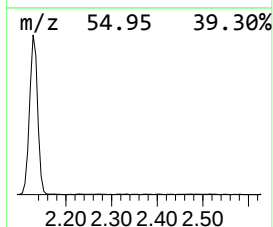
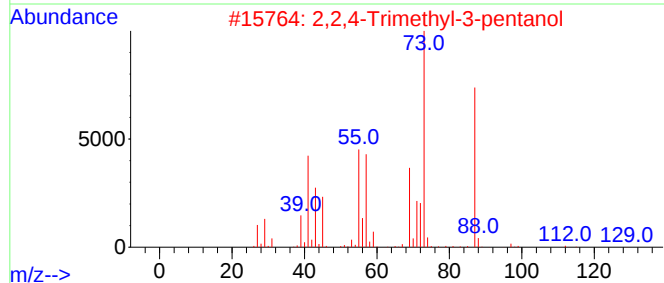
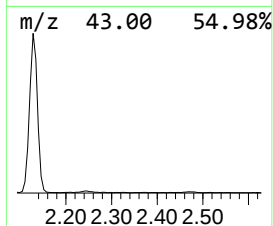
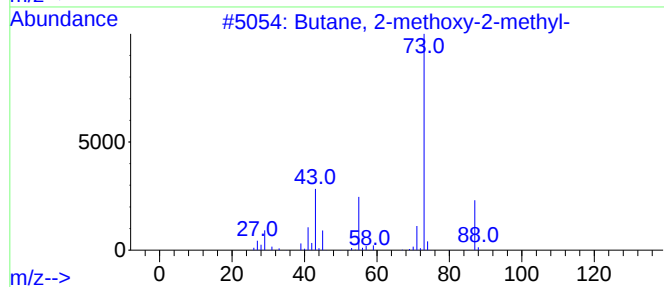
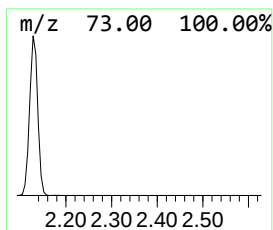
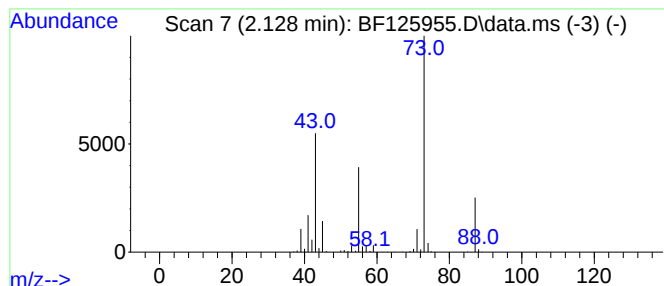
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.128	28.98 ng	1711480	1,4-Dichlorobenzene-d4	6.869

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Butane, 2-methoxy-2-methyl-	102	C6H14O	000994-05-8	83
2			2,2,4-Trimethyl-3-pentanol	130	C8H18O	005162-48-1	39
3			Silane, tetramethyl-	88	C4H12Si	000075-76-3	35
4			3-Pentanol, 3-methyl-	102	C6H14O	000077-74-7	23
5			1,3-Dioxolane, 2-propyl-	116	C6H12O2	003390-13-4	17



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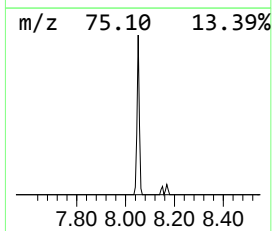
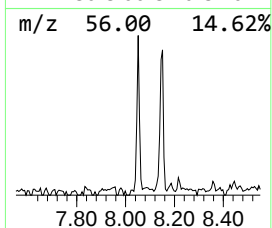
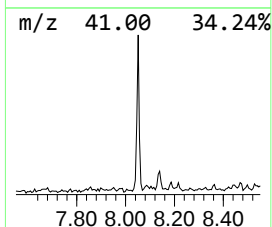
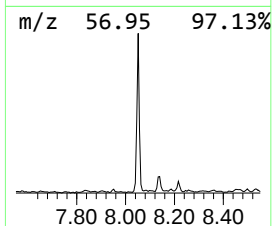
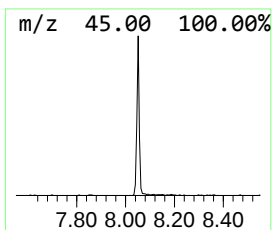
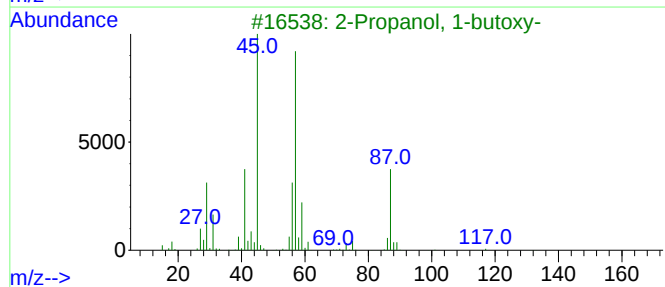
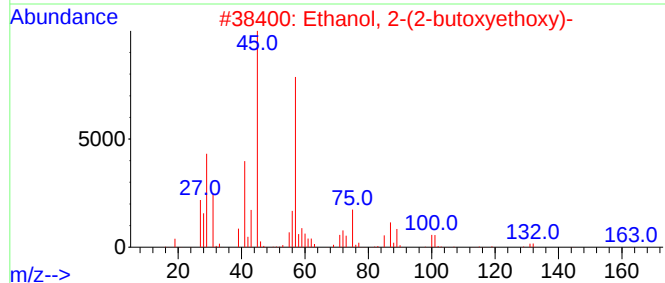
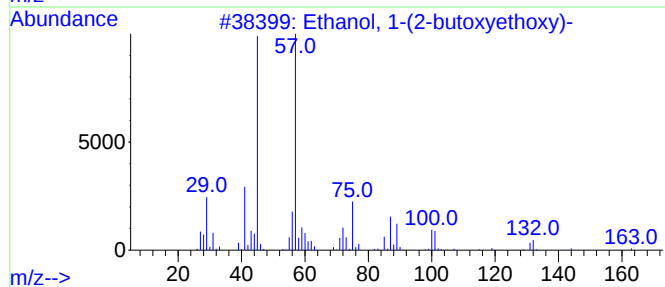
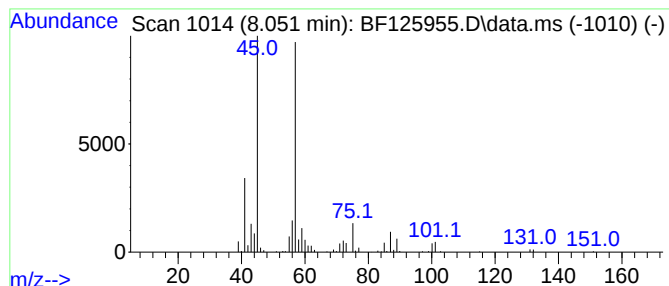
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 Ethanol, 1-(2-butoxyethoxy)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.051	3.71 ng	265923	Naphthalene-d8	8.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 1-(2-butoxyethoxy)-	162	C8H18O3	054446-78-5	78
2			Ethanol, 2-(2-butoxyethoxy)-	162	C8H18O3	000112-34-5	72
3			2-Propanol, 1-butoxy-	132	C7H16O2	005131-66-8	59
4			Ethylene glycol monoisobutyl ether	118	C6H14O2	004439-24-1	59
5			Ethanol, 2-[2-(2-methylpropoxy)e...	162	C8H18O3	018912-80-6	56



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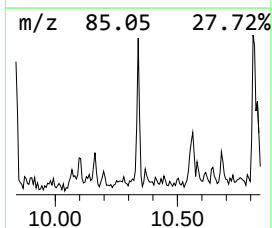
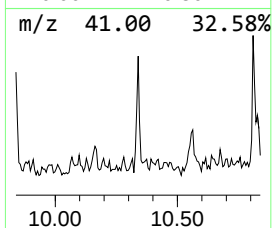
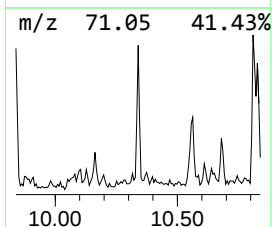
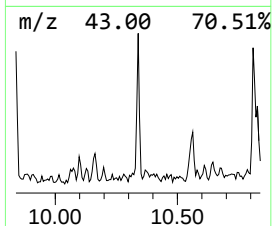
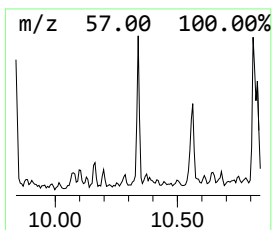
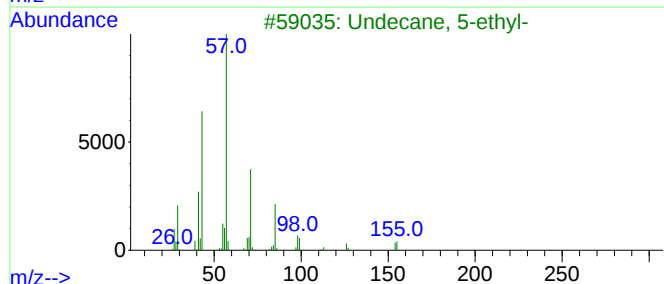
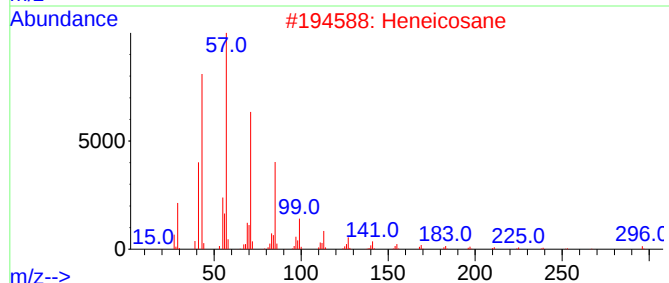
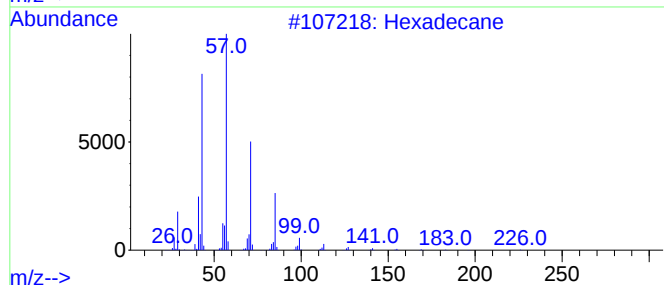
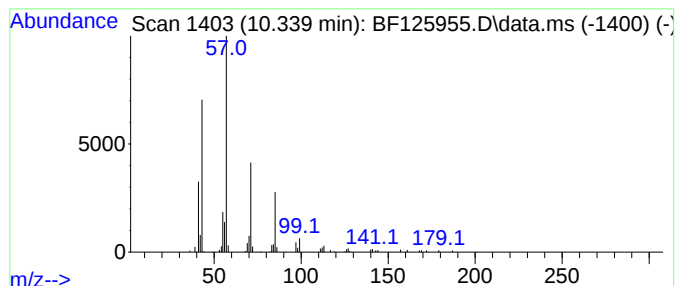
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TIC Library : C:\Database\NIST20.L
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Peak Number 3 Hexadecane Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.339	2.05 ng	130334	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	87
2			Heneicosane	296	C21H44	000629-94-7	78
3			Undecane, 5-ethyl-	184	C13H28	017453-94-0	78
4			Octacosane	394	C28H58	000630-02-4	78
5			Eicosane	282	C20H42	000112-95-8	78



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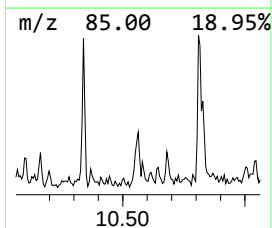
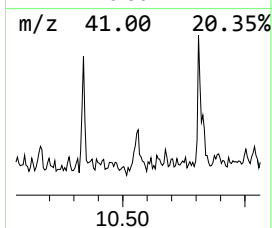
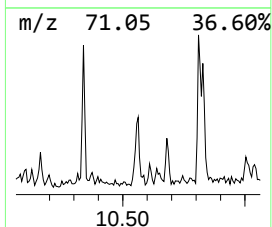
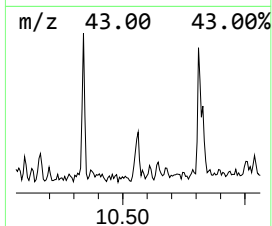
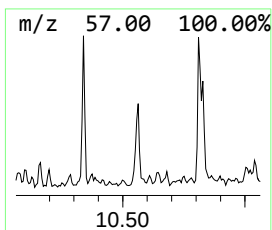
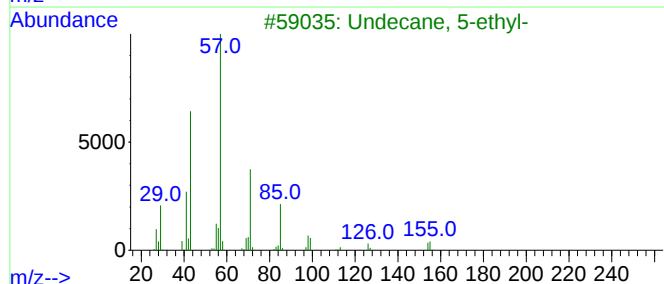
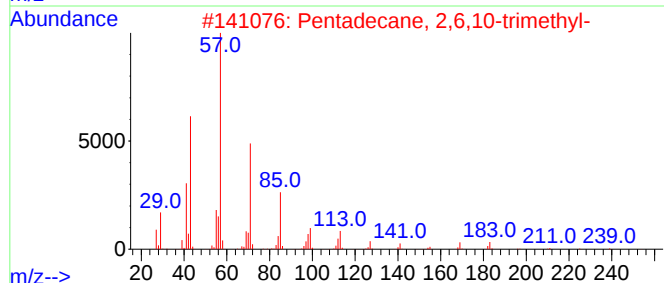
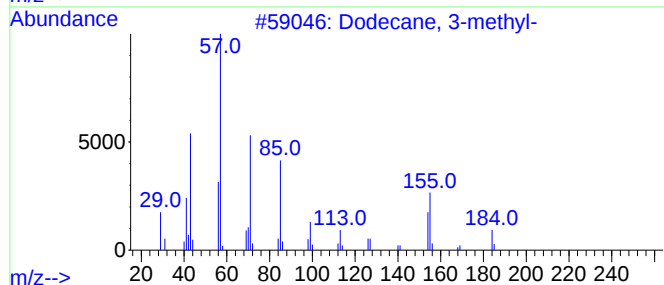
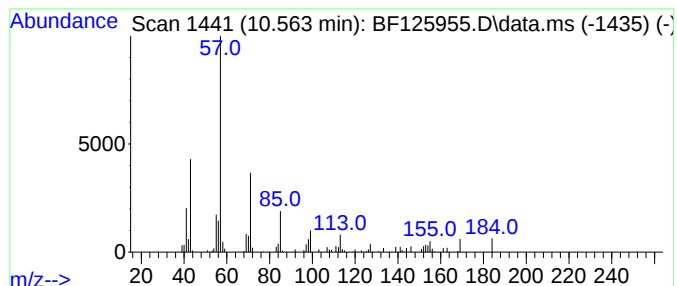
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TIC Library : C:\Database\NIST20.L
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Peak Number 4 Dodecane, 3-methyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.563	2.05 ng	129837	Acenaphthene-d10	9.904

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 3-methyl-	184	C13H28	017312-57-1	64
2			Pentadecane, 2,6,10-trimethyl-	254	C18H38	003892-00-0	64
3			Undecane, 5-ethyl-	184	C13H28	017453-94-0	59
4			Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	58
5			2-Bromo dodecane	248	C12H25Br	013187-99-0	53



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Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
Operator : JU/CG
Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
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ClientSampleId :
OK-2-102521

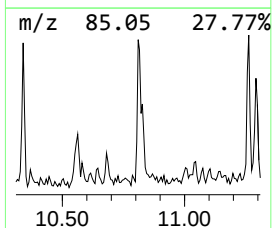
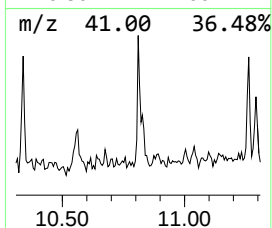
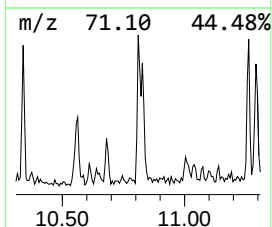
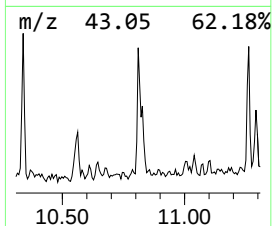
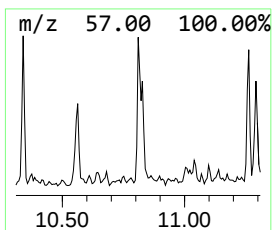
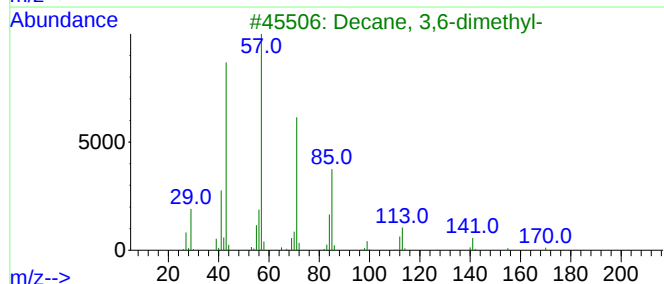
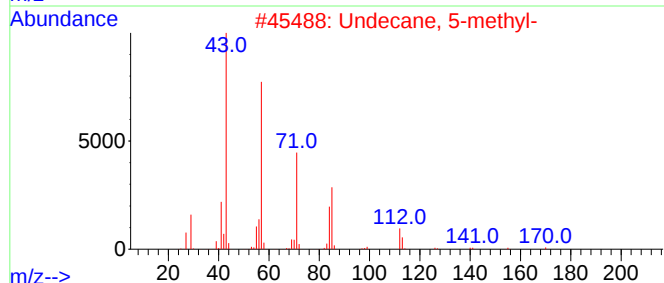
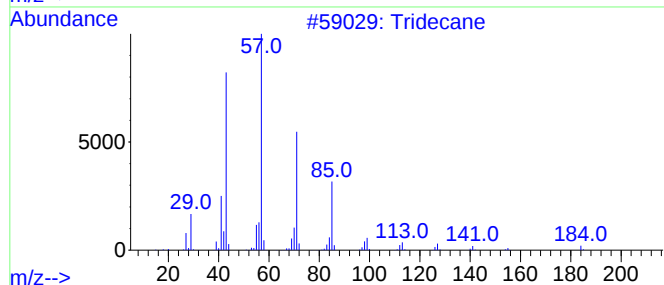
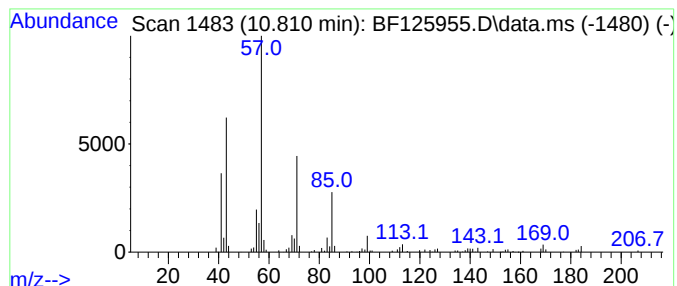
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 5 Tridecane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.810	3.98 ng	245993	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	90
2			Undecane, 5-methyl-	170	C12H26	001632-70-8	80
3			Decane, 3,6-dimethyl-	170	C12H26	017312-53-7	80
4			Dodecane, 1-iodo-	296	C12H25I	004292-19-7	72
5			5-Ethyldecane	170	C12H26	017302-36-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
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Sample : M4355-01 2X
Misc :
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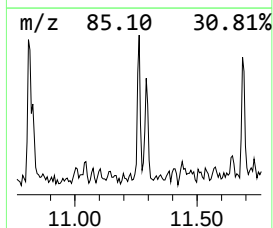
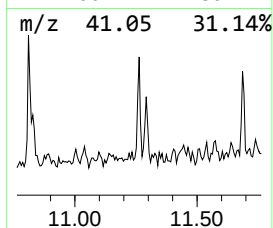
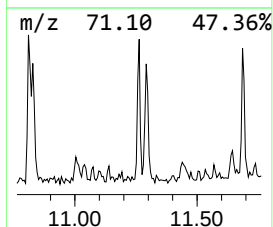
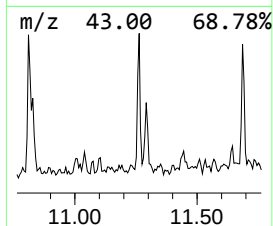
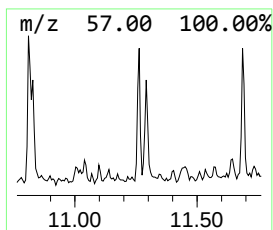
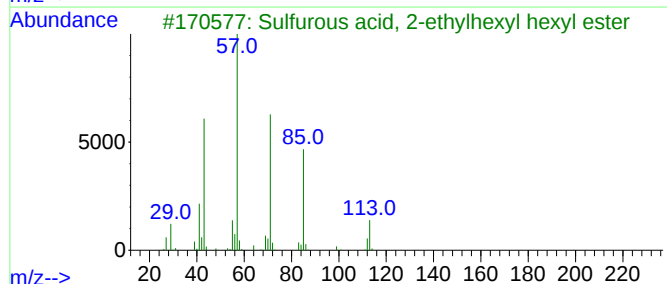
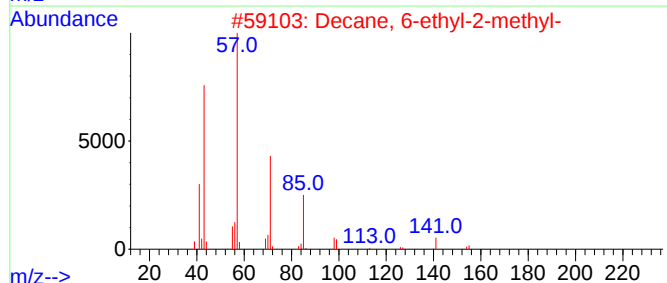
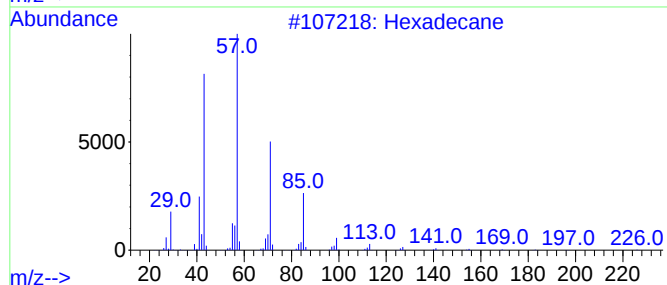
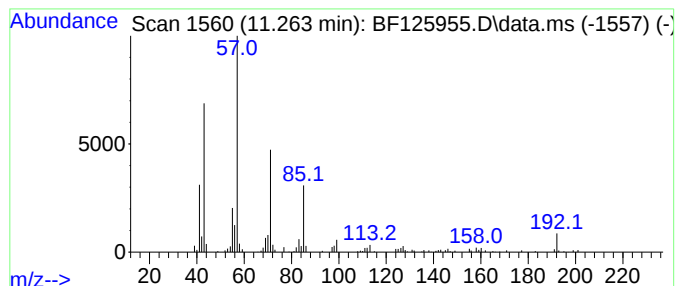
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ClientSampleId :
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 6 Decane, 6-ethyl-2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
11.263	2.06 ng	127025	Phenanthrene-d10		11.386	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	80
2	Decane, 6-ethyl-2-methyl-		184	C13H28	062108-21-8	72
3	Sulfurous acid, 2-ethylhexyl hex...		278	C14H30O3S	1000309-20-2	72
4	Decane, 3-methyl-		156	C11H24	013151-34-3	70
5	Nonadecane		268	C19H40	000629-92-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
Operator : JU/CG
Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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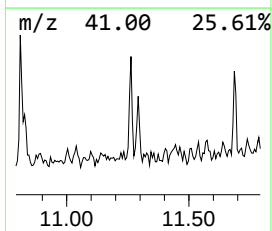
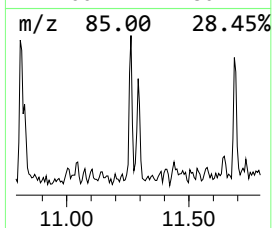
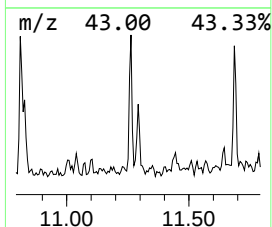
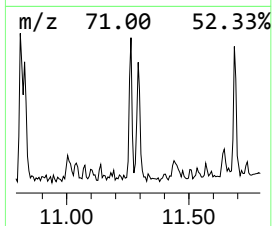
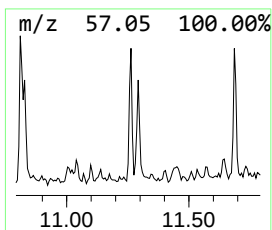
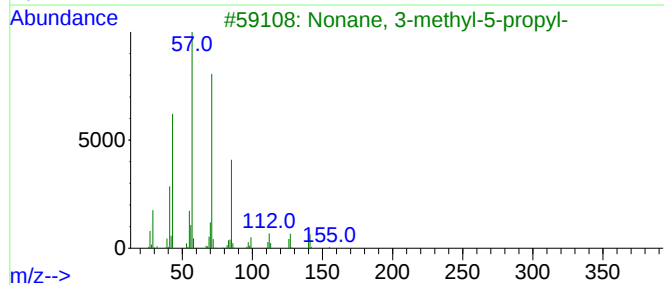
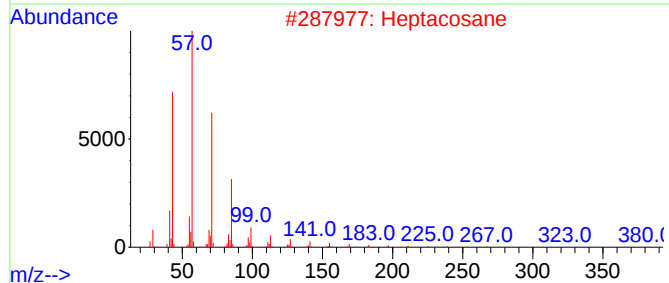
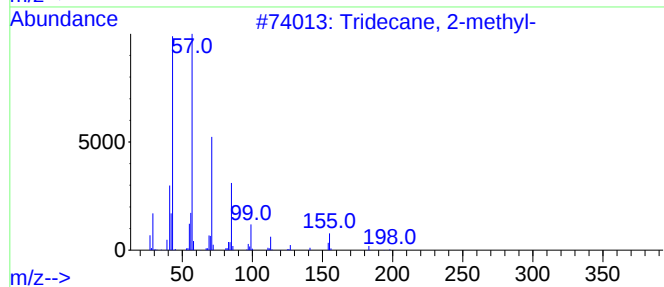
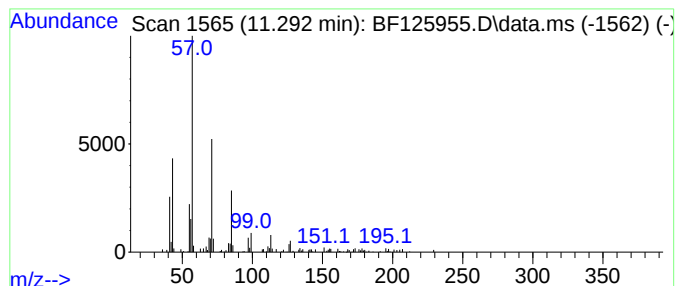
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 7 Tridecane, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.292	2.10 ng	129553	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 2-methyl-	198	C14H30	001560-96-9	83
2			Heptacosane	380	C27H56	000593-49-7	80
3			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	72
4			Tetradecane, 1-iodo-	324	C14H29I	019218-94-1	72
5			Undecane	156	C11H24	001120-21-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
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Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

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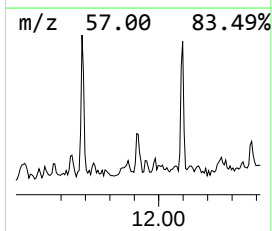
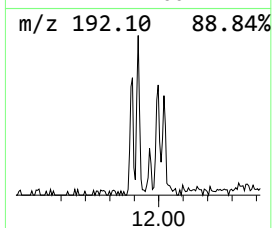
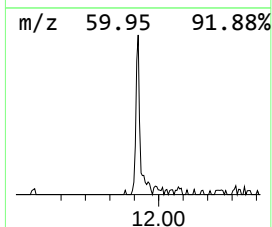
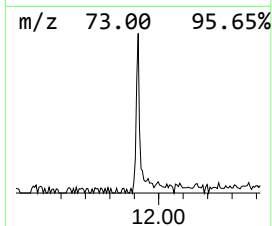
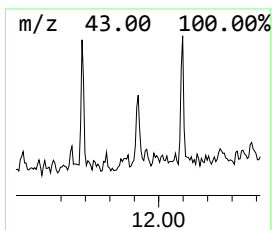
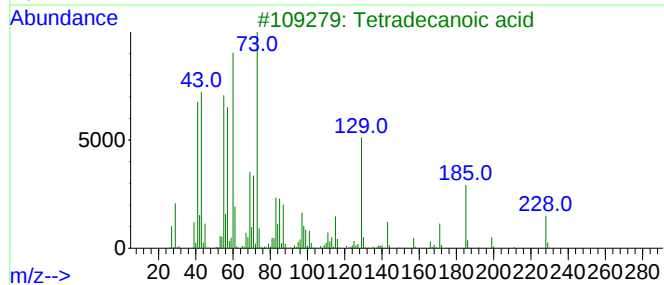
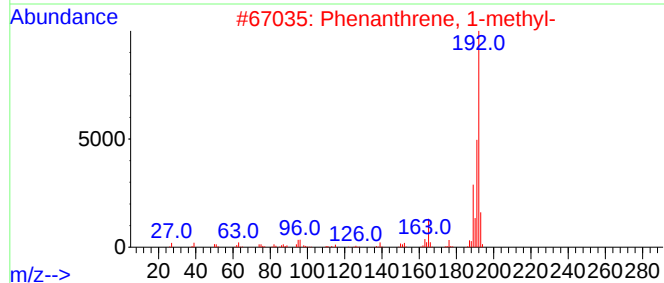
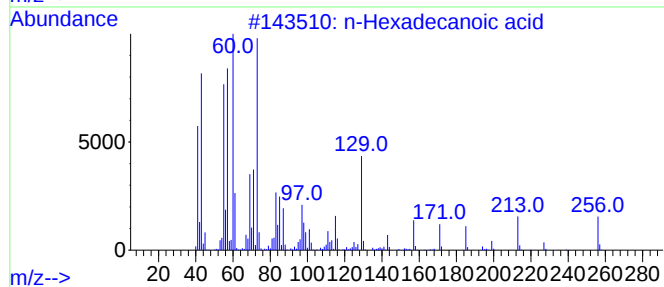
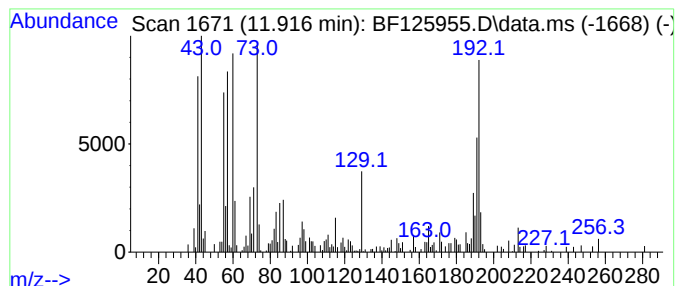
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.916	2.76 ng	170772	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	90
2			Phenanthrene, 1-methyl-	192	C15H12	000832-69-9	70
3			Tetradecanoic acid	228	C14H28O2	000544-63-8	60
4			Anthracene, 2-methyl-	192	C15H12	000613-12-7	50
5			Tridecanoic acid	214	C13H26O2	000638-53-9	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
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Misc :
ALS Vial : 22 Sample Multiplier: 1

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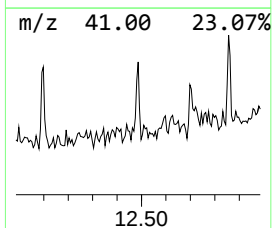
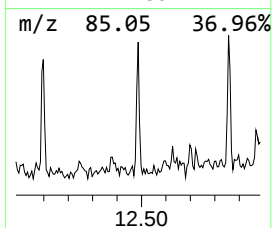
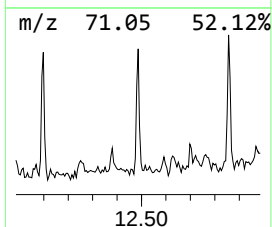
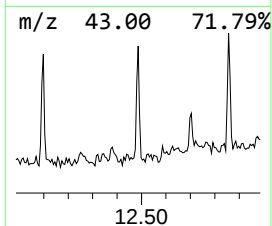
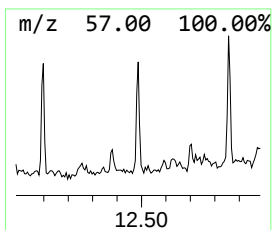
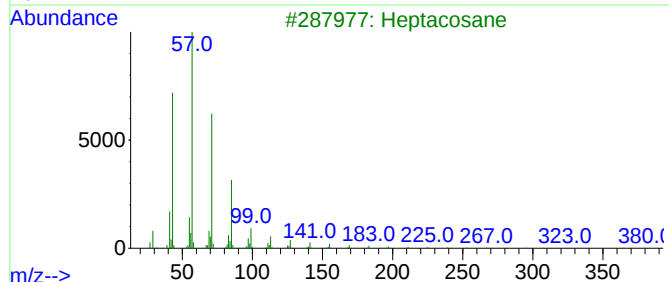
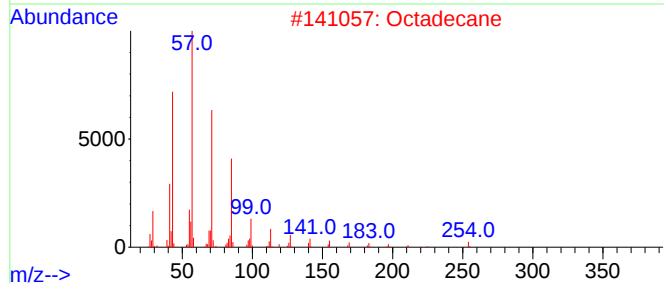
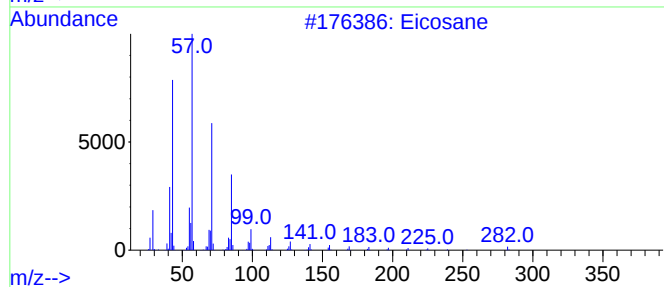
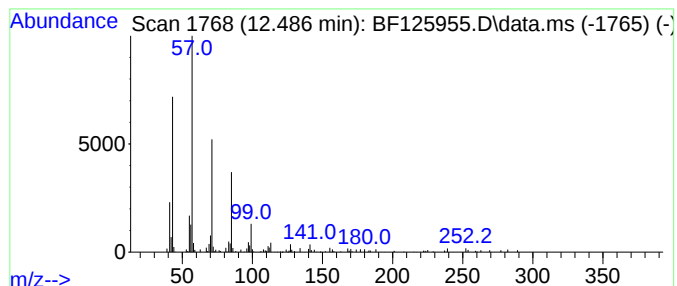
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 9 Eicosane Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.486	2.50 ng	154720	Phenanthrene-d10	11.386

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane	282	C20H42	000112-95-8	93
2			Octadecane	254	C18H38	000593-45-3	90
3			Heptacosane	380	C27H56	000593-49-7	83
4			2-Bromo dodecane	248	C12H25Br	013187-99-0	83
5			Tetracosane	338	C24H50	000646-31-1	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
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Sample : M4355-01 2X
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ClientSampleId :
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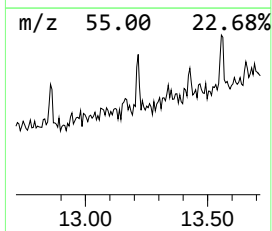
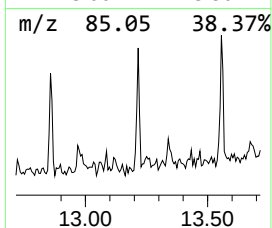
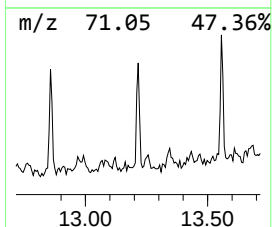
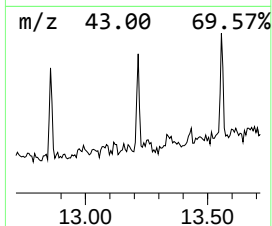
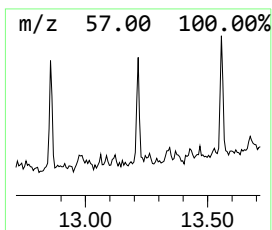
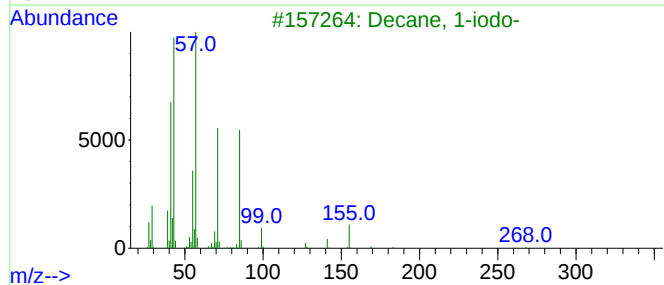
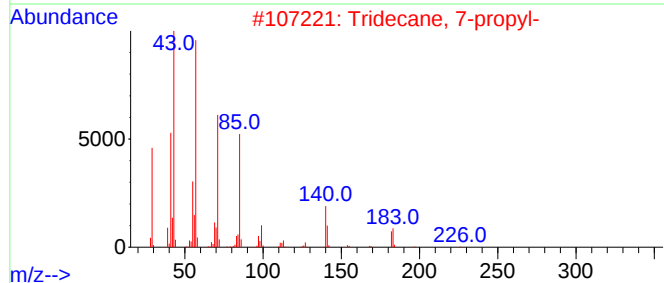
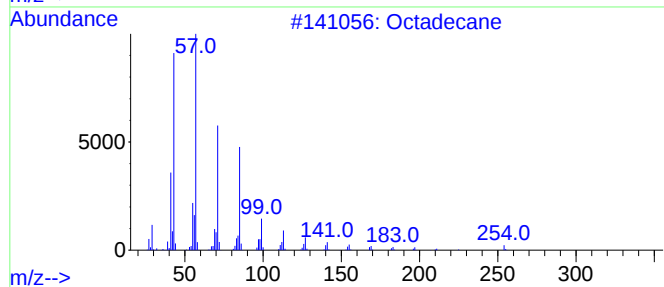
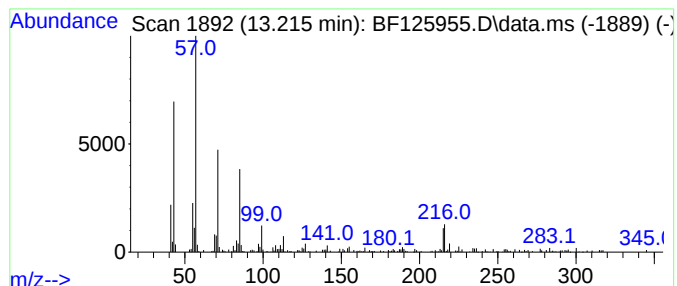
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 10 Tridecane, 7-propyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.216	3.80 ng	173142	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane	254	C18H38	000593-45-3	92
2			Tridecane, 7-propyl-	226	C16H34	055045-09-5	86
3			Decane, 1-iodo-	268	C10H21I	002050-77-3	86
4			Eicosane	282	C20H42	000112-95-8	76
5			Tetradecane, 4-ethyl-	226	C16H34	055045-14-2	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
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ClientSampleId :
OK-2-102521

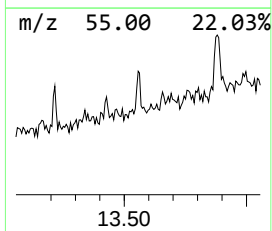
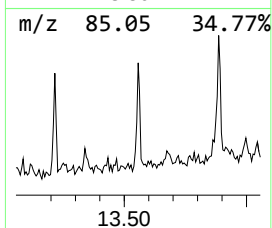
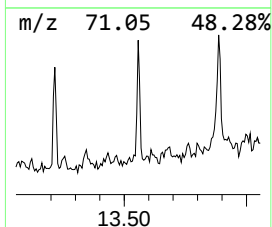
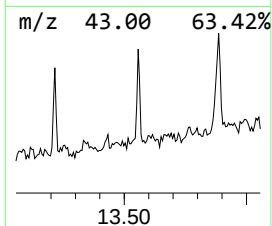
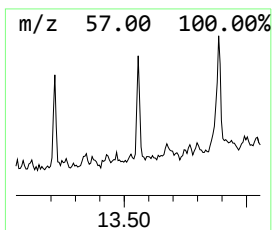
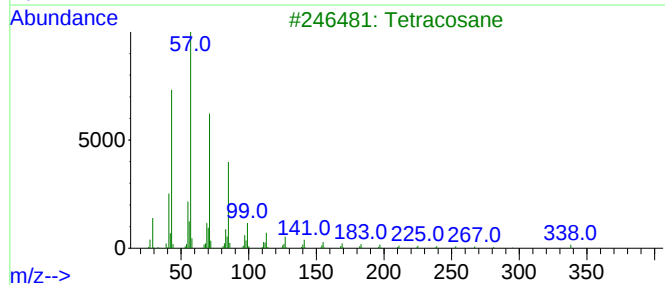
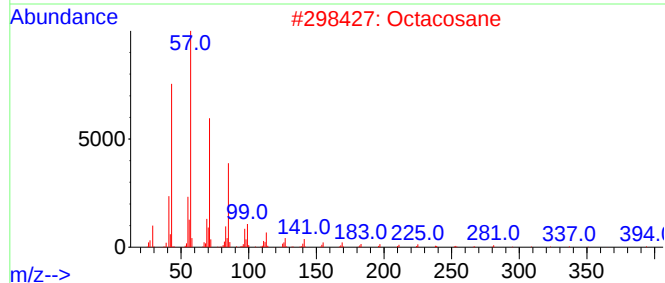
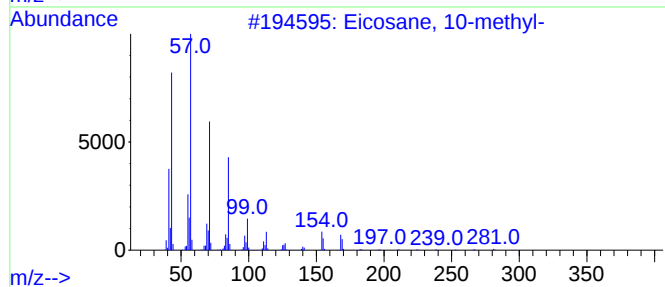
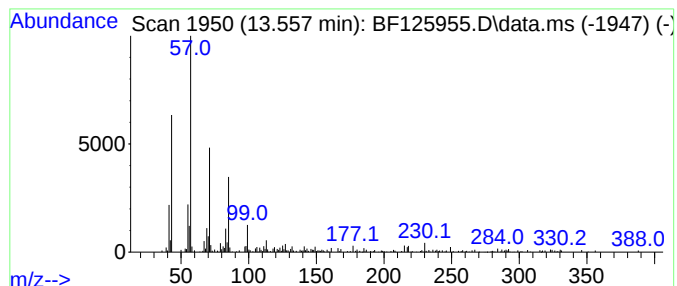
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 11 Eicosane, 10-methyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.557	4.46 ng	203043	Chrysene-d12	14.027

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Eicosane, 10-methyl-	296	C21H44	054833-23-7	76
2			Octacosane	394	C28H58	000630-02-4	74
3			Tetracosane	338	C24H50	000646-31-1	72
4			Heptacosane	380	C27H56	000593-49-7	72
5			Tridecane, 6-methyl-	198	C14H30	013287-21-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
Operator : JU/CG
Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102521

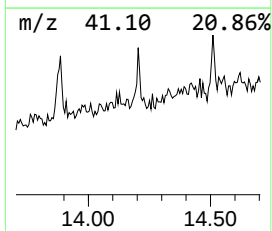
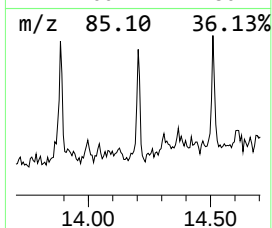
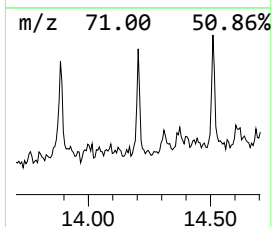
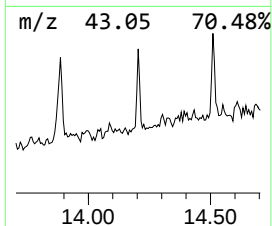
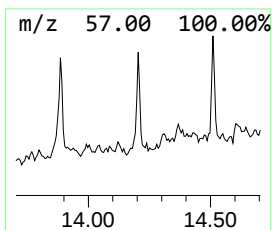
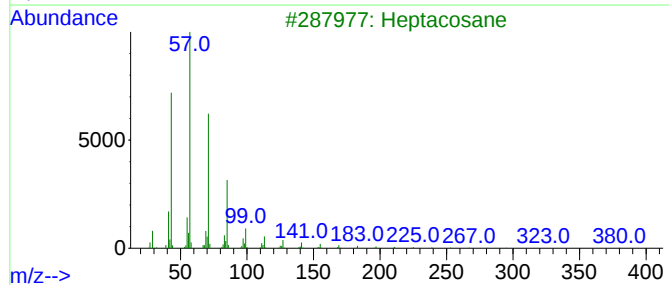
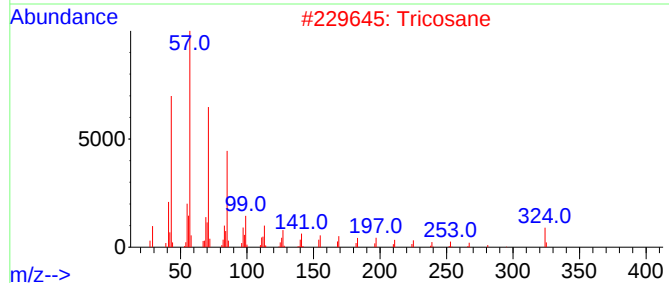
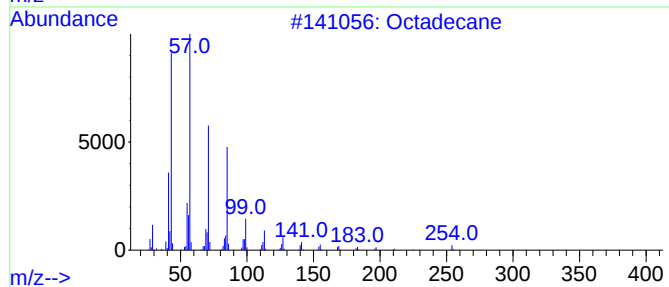
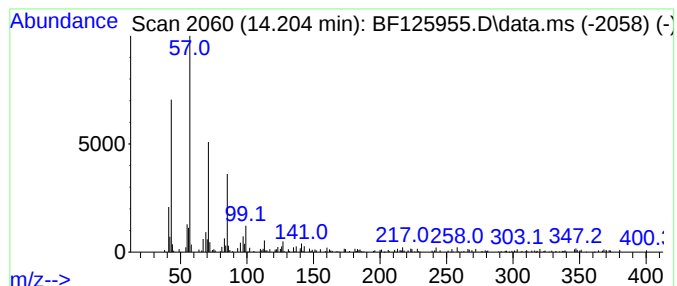
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 12 Octadecane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.204	3.83 ng	174419	Chrysene-d12	14.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	89
2		Tricosane	324	C23H48	000638-67-5	87
3		Heptacosane	380	C27H56	000593-49-7	83
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	80
5		Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
Operator : JU/CG
Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102521

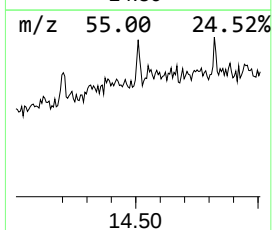
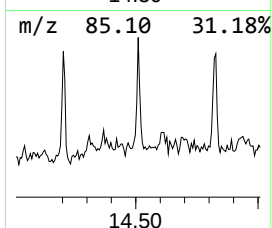
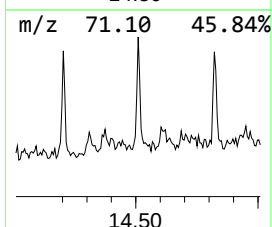
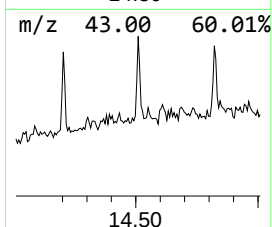
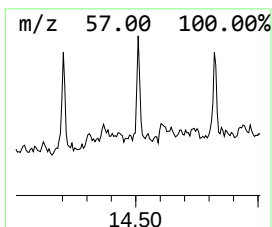
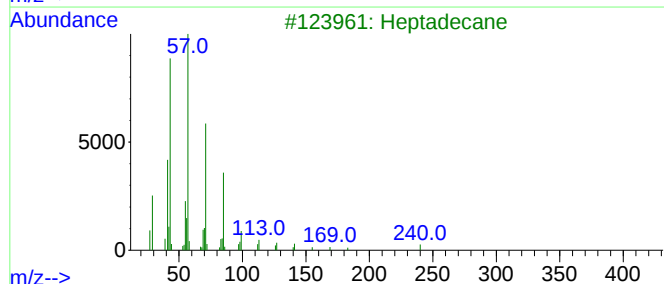
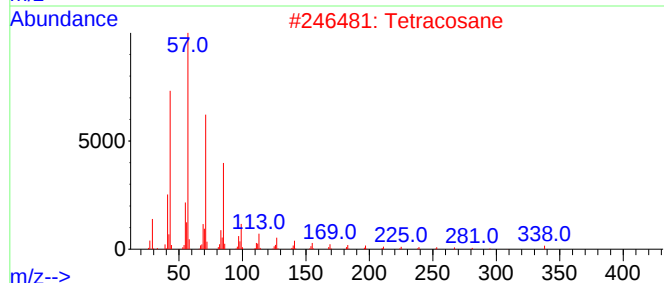
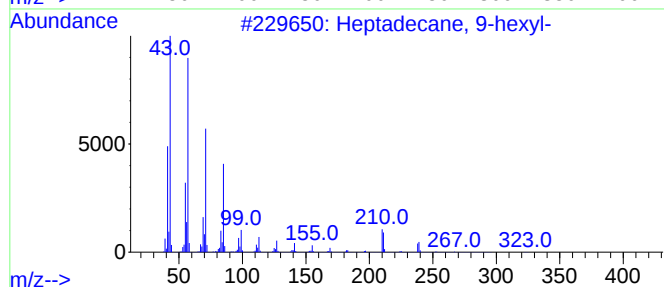
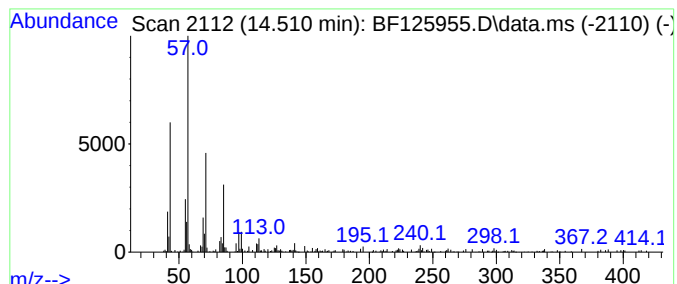
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

Peak Number 13 Heptadecane, 9-hexyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.509	4.17 ng	189842	Chrysene-d12	14.027

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	89
2		Tetracosane	338	C24H50	000646-31-1	89
3		Heptadecane	240	C17H36	000629-78-7	86
4		Hexadecane	226	C16H34	000544-76-3	86
5		Heptacosane, 1-chloro-	414	C27H55Cl	062016-79-9	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF102621\
Data File : BF125955.D
Acq On : 27 Oct 2021 2:30
Operator : JU/CG
Sample : M4355-01 2X
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
OK-2-102521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF102521.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...	2.128	29.0	ng	1711480	1	6.869	1181170	20.0
Ethanol, 1-(2-b...	8.051	3.7	ng	265923	2	8.151	1433600	20.0
Hexadecane	10.339	2.0	ng	130334	3	9.904	1268730	20.0
Dodecane, 3-met...	10.563	2.0	ng	129837	3	9.904	1268730	20.0
Tridecane	10.810	4.0	ng	245993	4	11.386	1236240	20.0
Decane, 6-ethyl...	11.263	2.1	ng	127025	4	11.386	1236240	20.0
Tridecane, 2-me...	11.292	2.1	ng	129553	4	11.386	1236240	20.0
n-Hexadecanoic ...	11.916	2.8	ng	170772	4	11.386	1236240	20.0
Eicosane	12.486	2.5	ng	154720	4	11.386	1236240	20.0
Tridecane, 7-pr...	13.216	3.8	ng	173142	5	14.027	911094	20.0
Eicosane, 10-me...	13.557	4.5	ng	203043	5	14.027	911094	20.0
Octadecane	14.204	3.8	ng	174419	5	14.027	911094	20.0
Heptadecane, 9-...	14.509	4.2	ng	189842	5	14.027	911094	20.0