LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020821\
Data File : BF123156.D
Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-020521

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-BF012721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF123156.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.181	9	16	22	rVB	507055	668892	13.05%	1.281%
2	5.510	577	582	585	rBV	3856778	3863894	75.36%	7.399%
3	6.516	748	753	756	rBV	3742161	3932200	76.69%	7.530%
4	6.892	813	817	820	rBV	1504772	1180093	23.02%	2.260%
5	7.451	904	912	915	rBV	2591956	2639522	51.48%	5.054%
6	7.892	977	987	989	rBV4	30831	69250	1.35%	0.133%
7	8.175	1029	1035	1044	rBV	1776983	1611799	31.44%	3.086%
8	8.769	1133	1136	1140	rBV	73209	57944	1.13%	0.111%
9	8.886	1154	1156	1160	rBV	162188	140071	2.73%	0.268%
10	8.986	1169	1173	1178	rBV	167280	153409	2.99%	0.294%
11	9.257	1214	1219	1222	rBV	4837836	4777085	93.17%	9.148%
12	9.333	1228	1232	1234	rBV2	86067	69629	1.36%	0.133%
13	9.516	1259	1263	1268	rBV2	67488	84782	1.65%	0.162%
14	9.592	1272	1276	1278	rVV	78620	69840	1.36%	0.134%
15	9.645	1282	1285	1289	rVV	212724	182952	3.57%	0.350%
16	9.863	1318	1322	1326	rVB	78041	66282	1.29%	0.127%
17	9.933	1327	1334	1337	rBV	2048789	1814030	35.38%	3.474%
18	9.969	1337	1340	1343	rVB	494626	424673	8.28%	0.813%
19	10.139	1365	1369	1373	rBV	191334	183666	3.58%	0.352%
20	10.363	1405	1407	1410	rVB	74492	67136	1.31%	0.129%
21	10.486	1424	1428	1433	rVB	293133	293235	5.72%	0.562%
22	10.727	1464	1469	1480	rBV	2904943	2869977	55.97%	5.496%
23	10.839	1486	1488	1493	rVB2	81901	99205	1.93%	0.190%
24	11.286	1558	1564	1567	rBV3	92973	136451	2.66%	0.261%
25	11.322	1567	1570	1575	rVB	126748	120645	2.35%	0.231%
26	11.427	1584	1588	1590	rBV	2108139	1845836	36.00%	3.535%
27	11.451	1590	1592	1595	rVV	1458408	1109556	21.64%	2.125%
28	11.498	1595	1600	1604	rVB	407626	447515	8.73%	0.857%
29	11.657	1621	1627	1635	rBV	171636	239526	4.67%	0.459%
30	11.722	1635	1638	1640	rBV	75298	65997	1.29%	0.126%
31	11.822	1651	1655	1657	rBV	292214	261808	5.11%	0.501%
32	11.927	1670	1673	1675	rBV	107600	100958	1.97%	0.193%
33	11.957	1675	1678	1684	rVB	636221	563838	11.00%	1.080%
34	12.045	1689	1693	1695	rBV	323561	303070	5.91%	0.580%
35	12.127	1705	1707	1711	rVB2	61503	54453	1.06%	0.104%
36	12.210	1715	1721	1722	rBV2	113719	110867	2.16%	0.212%

LSC Area Percent Report

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Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-020521

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-BF012721.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	12.227	1722	1724	1726	rVB	86678	59918	1.17%	0.115%
38	12.480	1764	1767	1769	rBV	66969	71112	1.39%	0.136%
39	12.510	1769	1772	1773	rVV	1060304	873376	17.03%	1.672%
40	12.527	1773	1775	1783	rVB	2136585	2026853	39.53%	3.881%
41	12.616	1787	1790	1792	rBV	471598	373786	7.29%	0.716%
42	12.645	1792	1795	1798	rBV	2115398	1716360	33.47%	3.287%
43	12.745	1808	1812	1819	rBV	934405	961098	18.74%	1.840%
44	12.798	1819	1821	1824	rVB	79583	73353	1.43%	0.140%
45	12.874	1828	1834	1837	rBV	1486799	1562352	30.47%	2.992%
46	12.921	1840	1842	1848	rVB2	69150	88065	1.72%	0.169%
47	13.021	1852	1859	1862	rBV	5552483	5127301	100.00%	9.818%
48	13.098	1867	1872	1875	rBV3	70512	125649	2.45%	0.241%
49	13.180	1883	1886	1887	rBV2	59509	61884	1.21%	0.119%
50	13.204	1887	1890	1894	rVB	330595	289935	5.65%	0.555%
51	13.251	1895	1898	1899	rBV2	62859	56820	1.11%	0.109%
52	13.268	1899	1901	1904	rVB	313712	253244	4.94%	0.485%
53	13.310	1904	1908	1910	rBV3	102948	106642	2.08%	0.204%
54	13.339	1910	1913	1916	rVB2	81336	104553	2.04%	0.200%
55	13.592	1954	1956	1960	rBV2	80834	88170	1.72%	0.169%
56	13.857	1998	2001	2003	rBV	72657	63147	1.23%	0.121%
57	13.880	2003	2005	2009	rVV2	69676	102898	2.01%	0.197%
58	13.921	2010	2012	2016	rVB	103485	97925	1.91%	0.188%
59	14.015	2021	2028	2032	rBV5	258272	632548	12.34%	1.211%
60	14.080	2034	2039	2041	rVV2	1705419	2251845	43.92%	4.312%
61	14.104	2041	2043	2050	rVB	688203	599920	11.70%	1.149%
62	14.186	2054	2057	2060	rVB	180014	152002	2.96%	0.291%
63	14.239	2064	2066	2069	rVB	98721	80389	1.57%	0.154%
64	14.551	2116	2119	2123	rVB3	109198	138199	2.70%	0.265%
65	14.862	2169	2172	2175	rBV	106657	104452	2.04%	0.200%
66	15.133	2214	2218	2221	rBV	445387	720974	14.06%	1.381%
67	15.439	2267	2270	2276	rVB	262746	339504	6.62%	0.650%
68	15.504	2277	2281	2287	rBV	365635	476769	9.30%	0.913%
69	15.568	2288	2292	2296	rBV	952244	1222252	23.84%	2.341%
70	17.068	2542	2547	2556	rVB2	154653	329268	6.42%	0.631%
71	17.521	2619	2624	2631	rBV	106491	208832	4.07%	0.400%

Sum of corrected areas: 52221481

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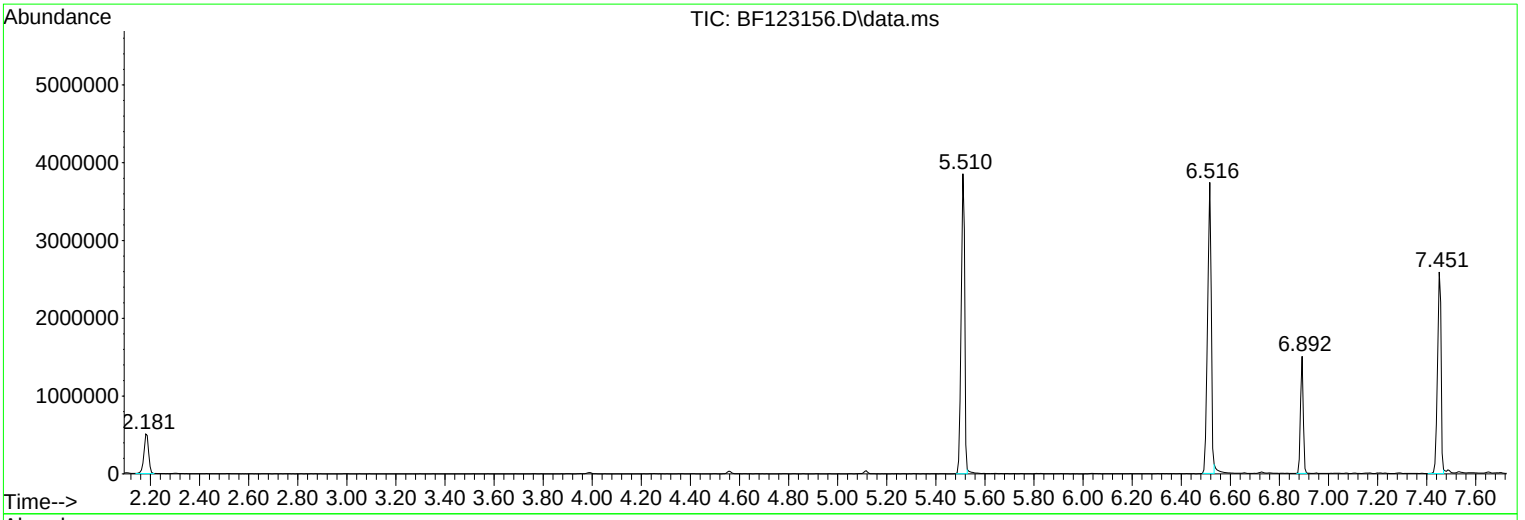
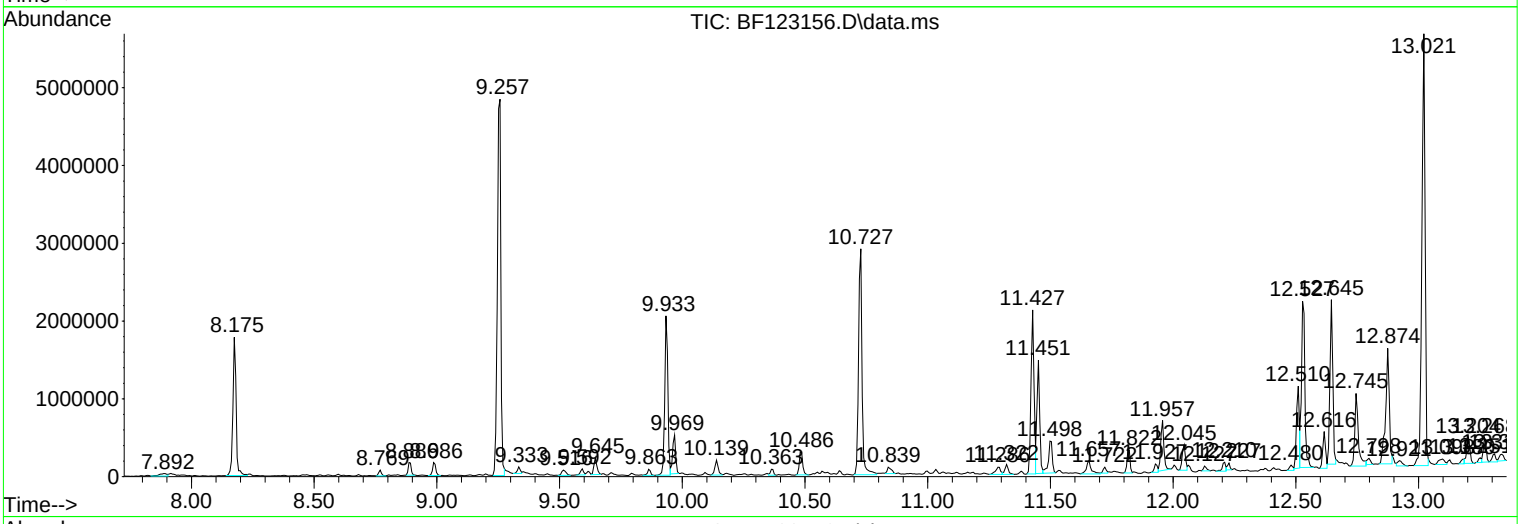
Misc :
ALS Vial : 18 Sample Multiplier: 1

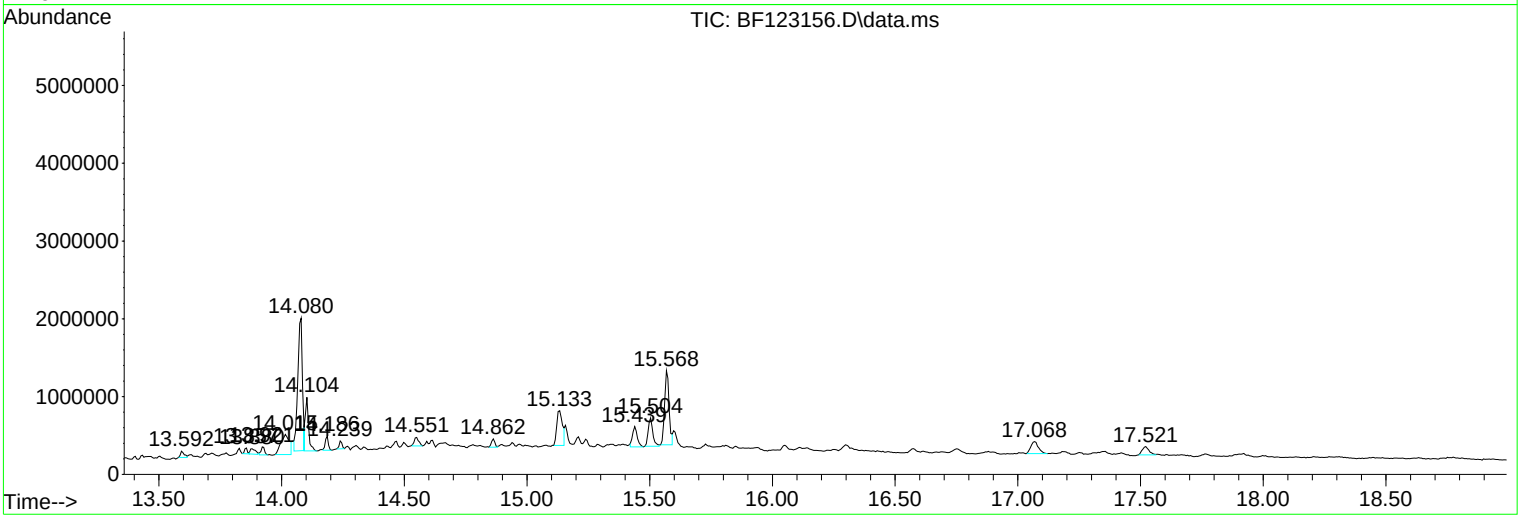
Instrument :
BNA_F
ClientSampleId :
TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P







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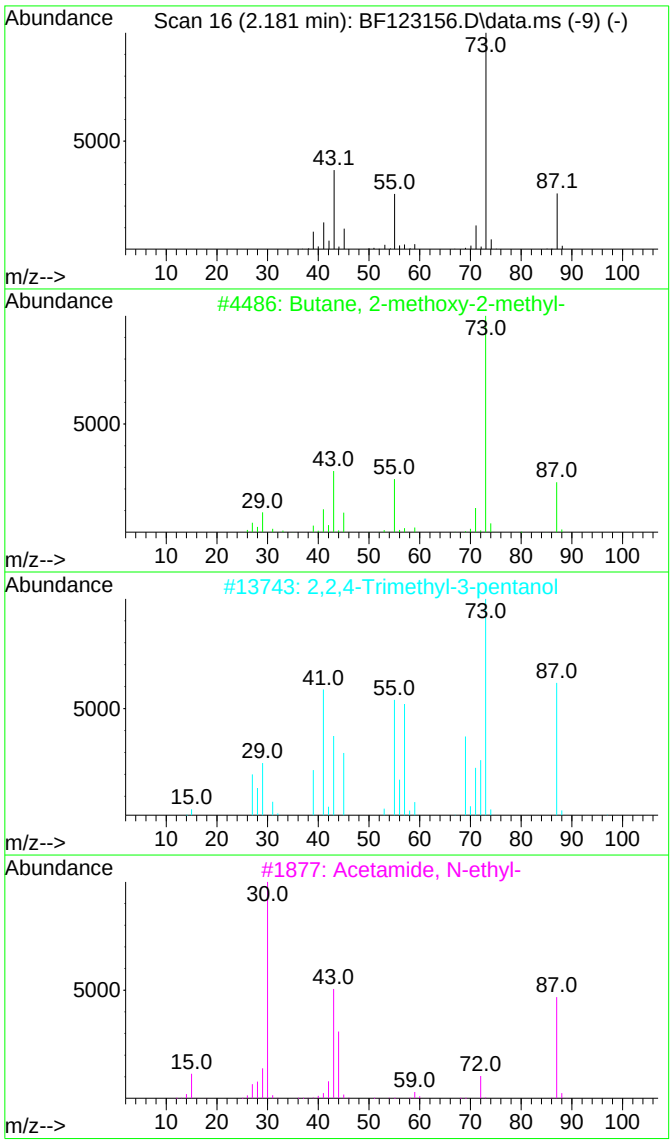
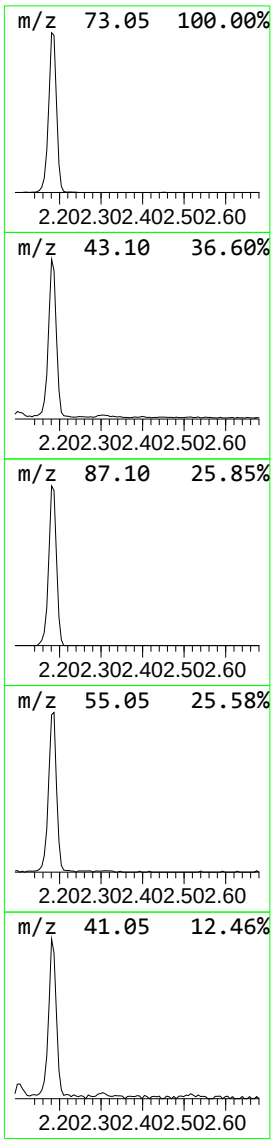
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.181	11.34 ng	668892	1,4-Dichlorobenzene-d4	6.892

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	40
3	Acetamide, N-ethyl-	87	C4H9NO	000625-50-3	22
4	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	17
5	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	9





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020821\
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Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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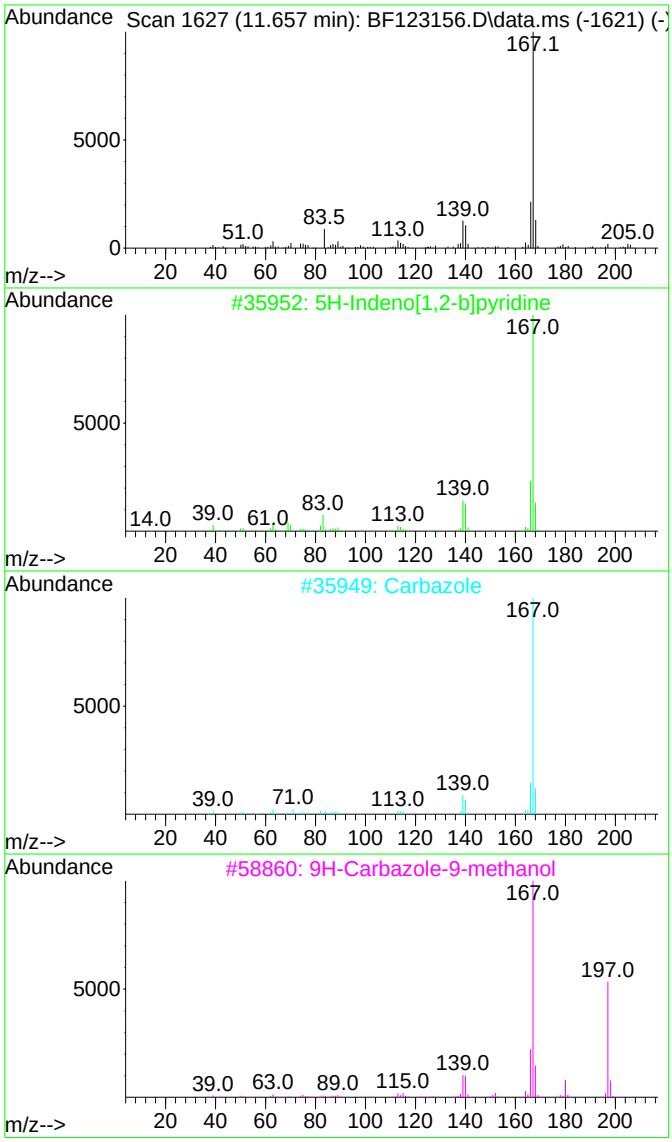
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

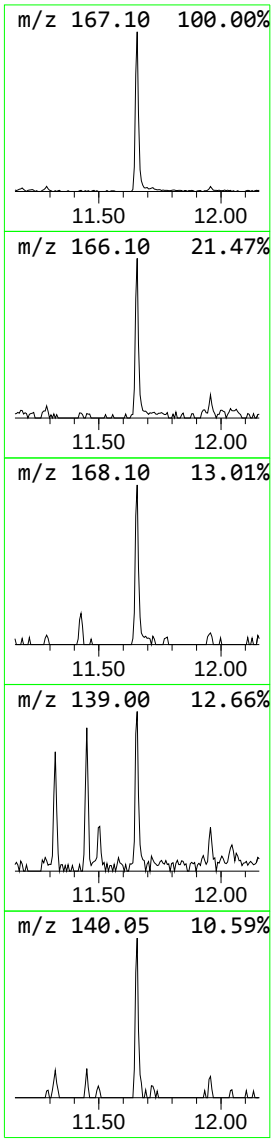
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 3 5H-Indeno[1,2-b]pyridine Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.657	2.60 ng	239526	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5H-Indeno[1,2-b]pyridine	167	C12H9N	000244-99-5	97
2	Carbazole	167	C12H9N	000086-74-8	91
3	9H-Carbazole-9-methanol	197	C13H11NO	002409-36-1	90
4	9H-Carbazole, 9-nitroso-	196	C12H8N2O	002788-23-0	90
5	D-Galactitol, 2-(acetylmethylami...	363	C16H29NO8	074410-42-7	90





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020821\
Data File : BF123156.D
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Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

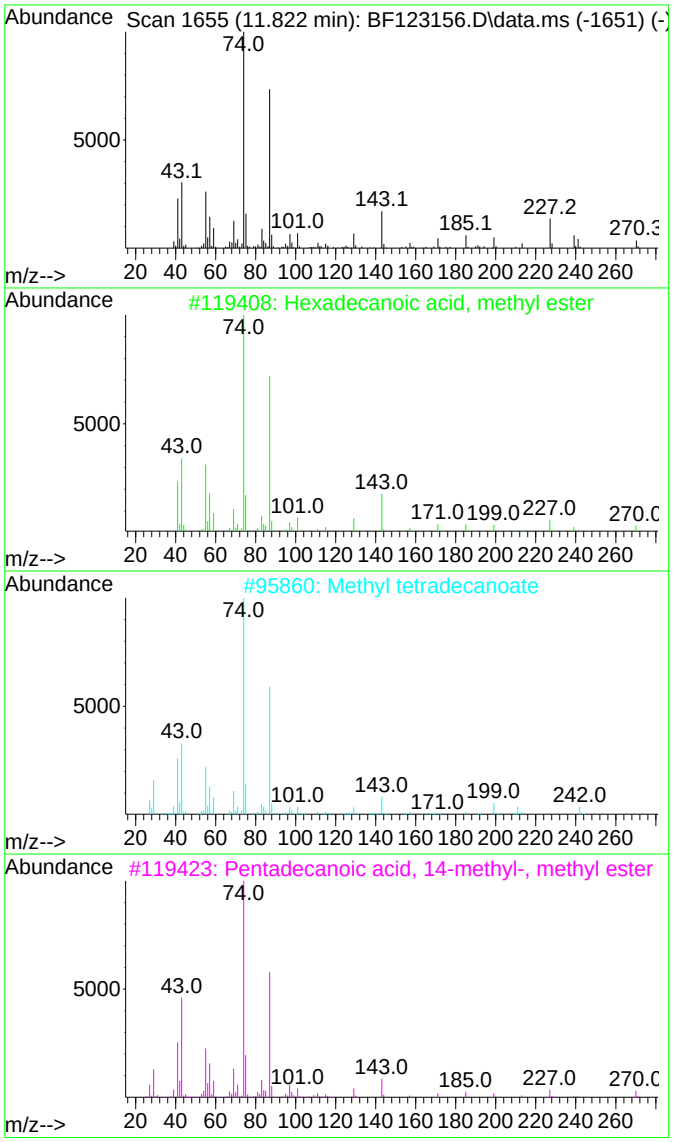
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

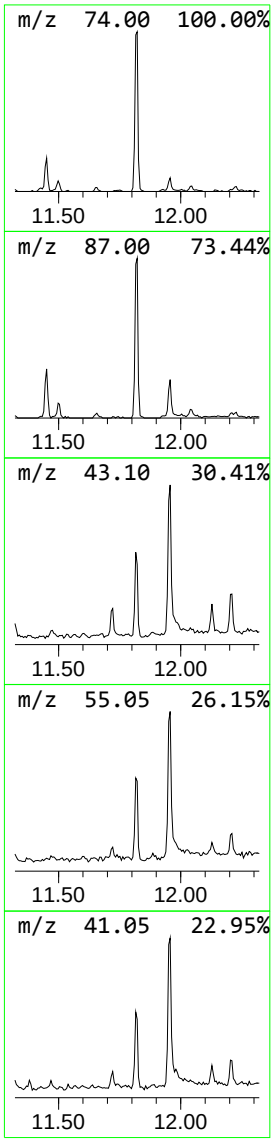
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 4 Hexadecanoic acid, methyl e... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.822	2.84 ng	261808	Phenanthrene-d10	11.427

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





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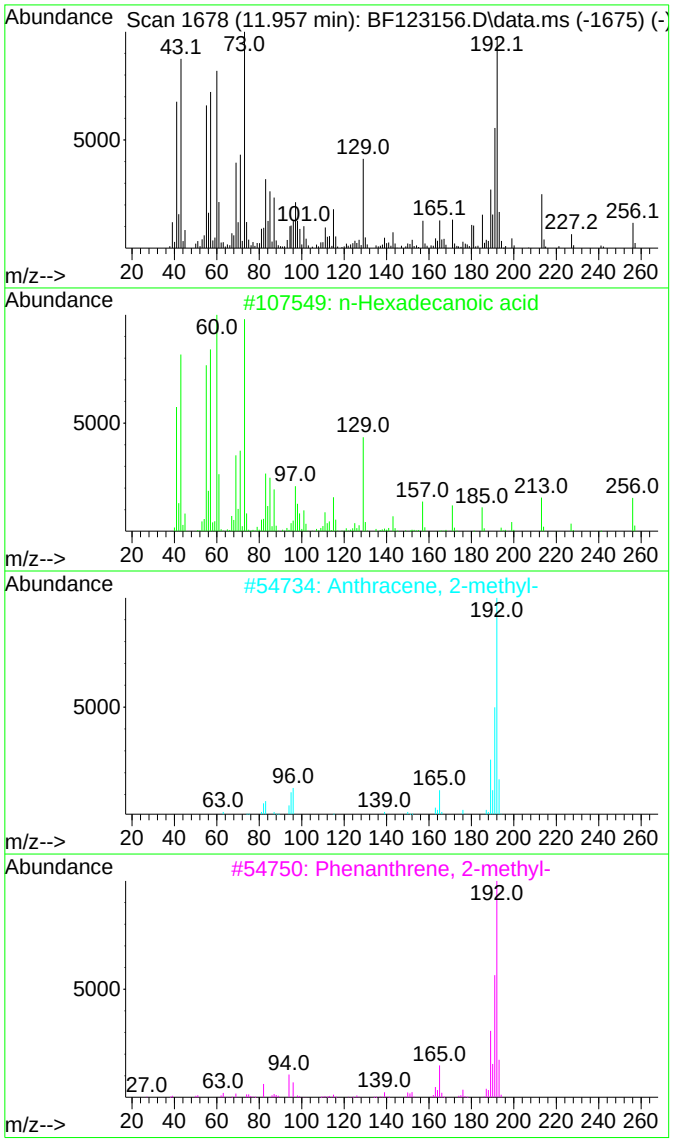
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

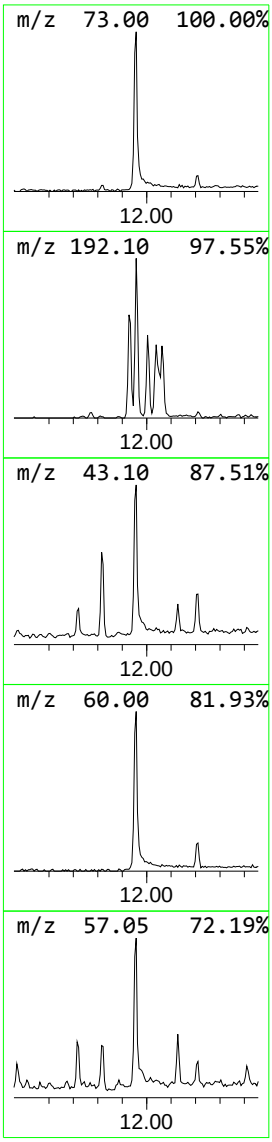
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 5 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.957	6.11 ng	563838	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
2	Anthracene, 2-methyl-	192	C15H12	000613-12-7	93
3	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	91
4	Anthracene, 9-methyl-	192	C15H12	000779-02-2	91
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	91





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Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

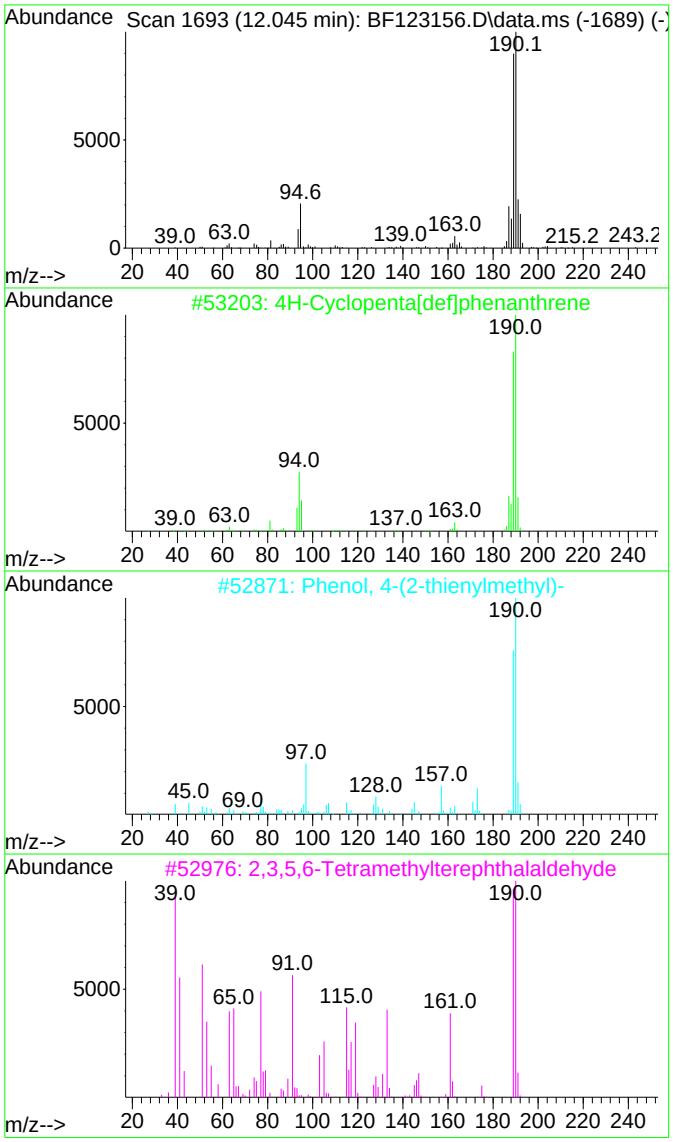
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

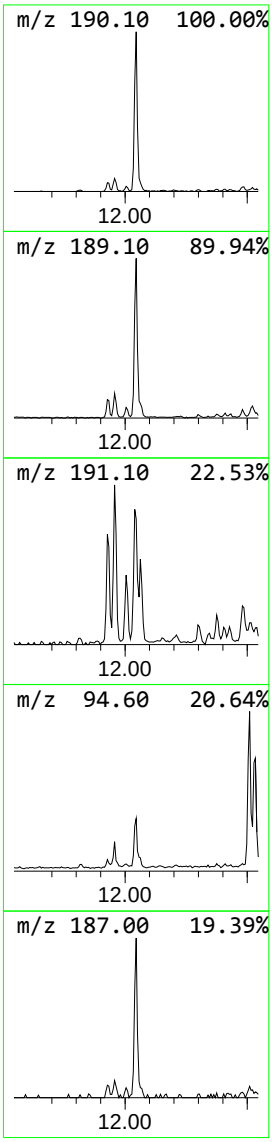
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 6 4H-Cyclopenta[def]phenanthrene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.045	3.28 ng	303070	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Cyclopenta[def]phenanthrene	190	C15H10	000203-64-5	90
2	Phenol, 4-(2-thienylmethyl)-	190	C11H10OS	091680-55-6	59
3	2,3,5,6-Tetramethylterephthalald...	190	C12H14O2	007072-01-7	50
4	2,2'-Bis(4,5-dimethylimidazole)	190	C10H14N4	069286-06-2	45
5	Thiophene-3-carboxaldehyde, 5-ch...	190	C5H4BClO3S	036155-87-0	40





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Data File : BF123156.D
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Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

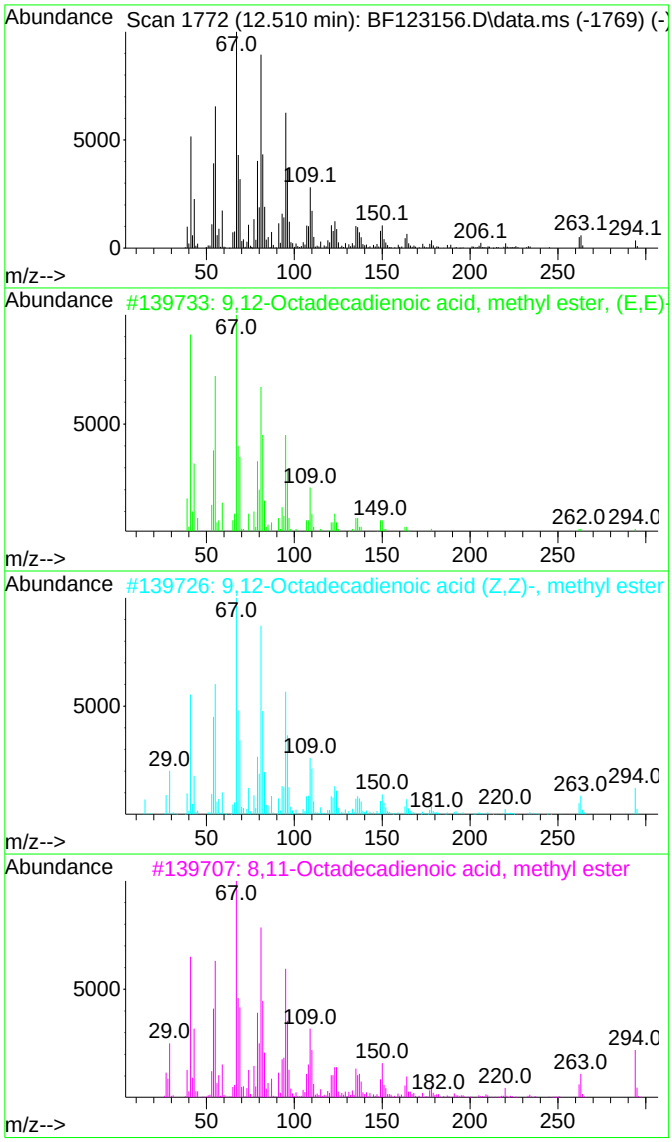
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

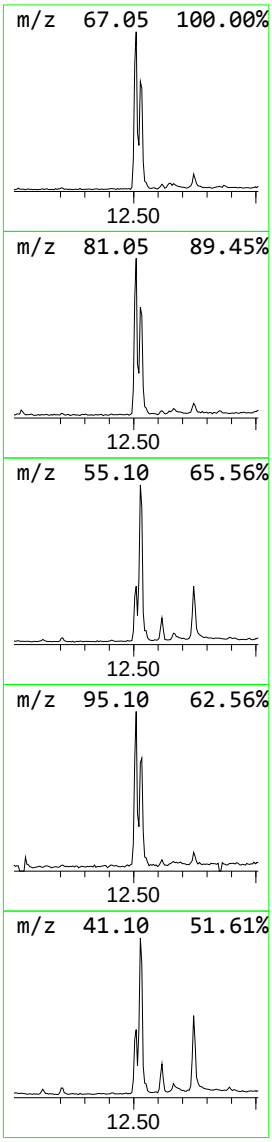
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 7 9,12-Octadecadienoic acid, ... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.510	9.46 ng	873376	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002566-97-4	99
2	9,12-Octadecadienoic acid (Z,Z)-...	294	C19H34O2	000112-63-0	99
3	8,11-Octadecadienoic acid, methy...	294	C19H34O2	056599-58-7	99
4	9,15-Octadecadienoic acid, methy...	294	C19H34O2	017309-05-6	99
5	9,12-Octadecadienoic acid, methy...	294	C19H34O2	002462-85-3	99





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020821\
Data File : BF123156.D
Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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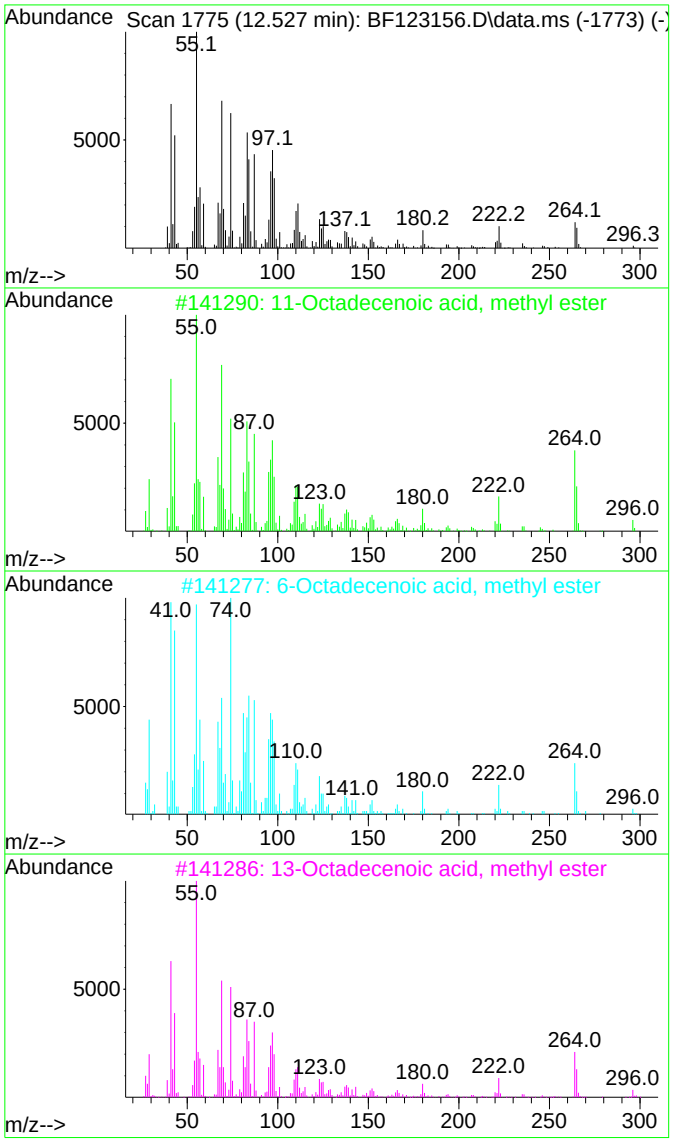
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

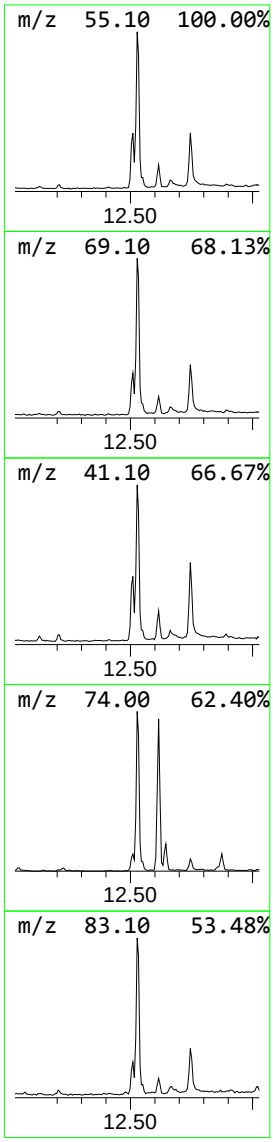
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 8 11-Octadecenoic acid, methy... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.527	21.96 ng	2026850	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	95
2	6-Octadecenoic acid, methyl ester	296	C19H36O2	052355-31-4	94
3	13-Octadecenoic acid, methyl ester	296	C19H36O2	056554-47-3	91
4	14-Octadecenoic acid, methyl ester	296	C19H36O2	056554-48-4	91
5	9-Octadecenoic acid (Z)-, methyl...	296	C19H36O2	000112-62-9	91





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Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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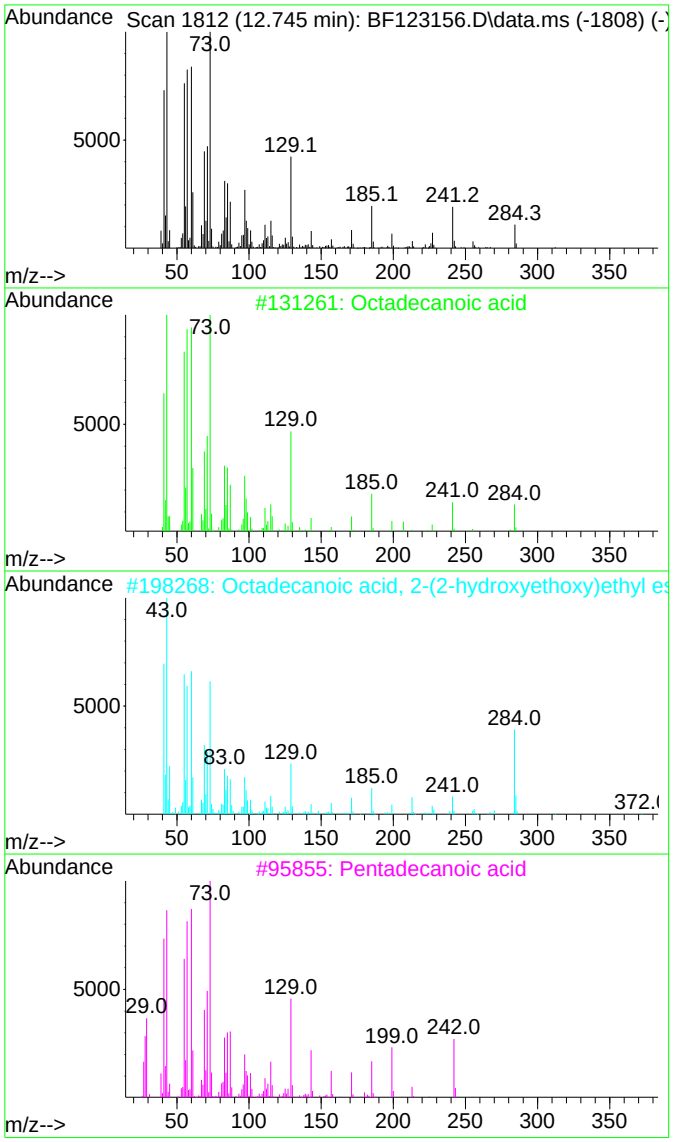
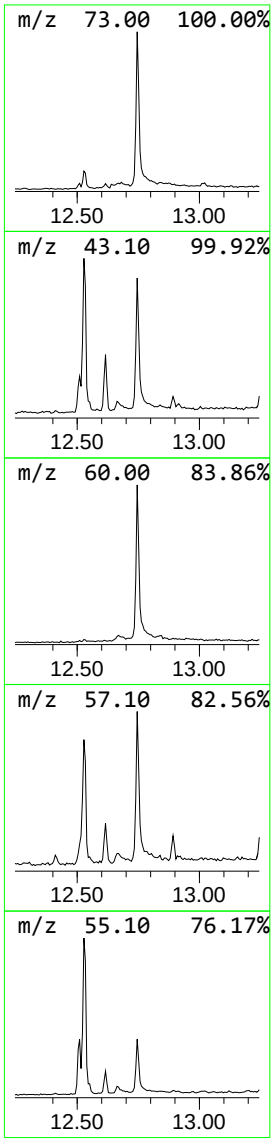
Peak Number 9 Octadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.745	10.41 ng	961098	Phenanthrene-d10	11.427

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecanoic acid	284	C18H36O2	000057-11-4	99
2	Octadecanoic acid, 2-(2-hydroxye...	372	C22H44O4	000106-11-6	91
3	Pentadecanoic acid	242	C15H30O2	001002-84-2	90
4	Tetradecanoic acid	228	C14H28O2	000544-63-8	86
5	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	74





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Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

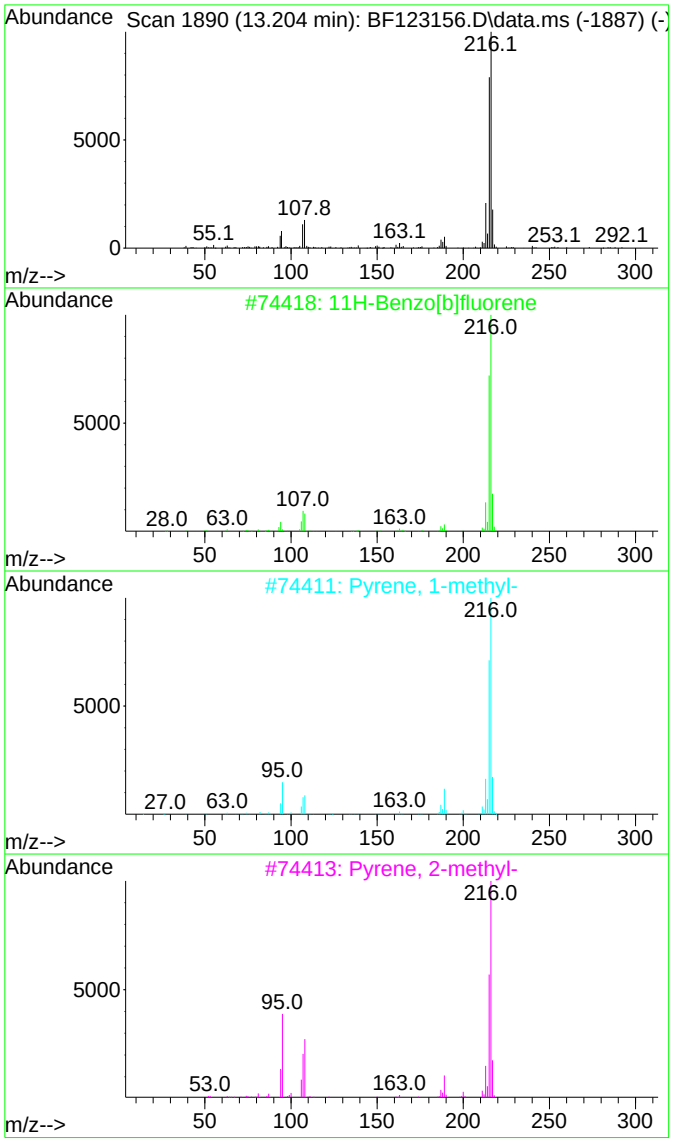
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

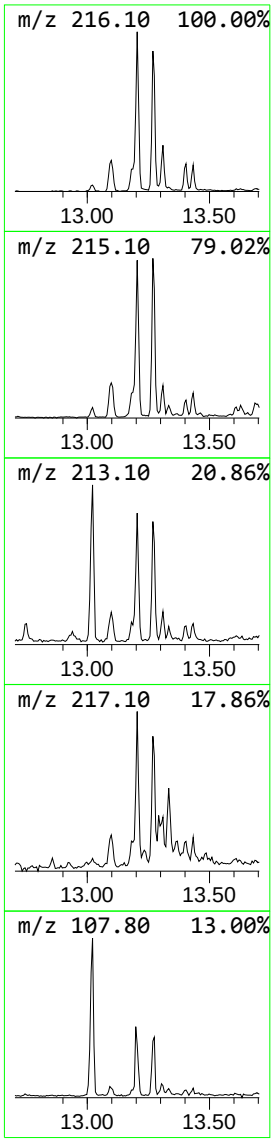
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 10 11H-Benzo[b]fluorene Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.204	2.58 ng	289935	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	94
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	93
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	87





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Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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TIC Integration Parameters: LSCINT.P

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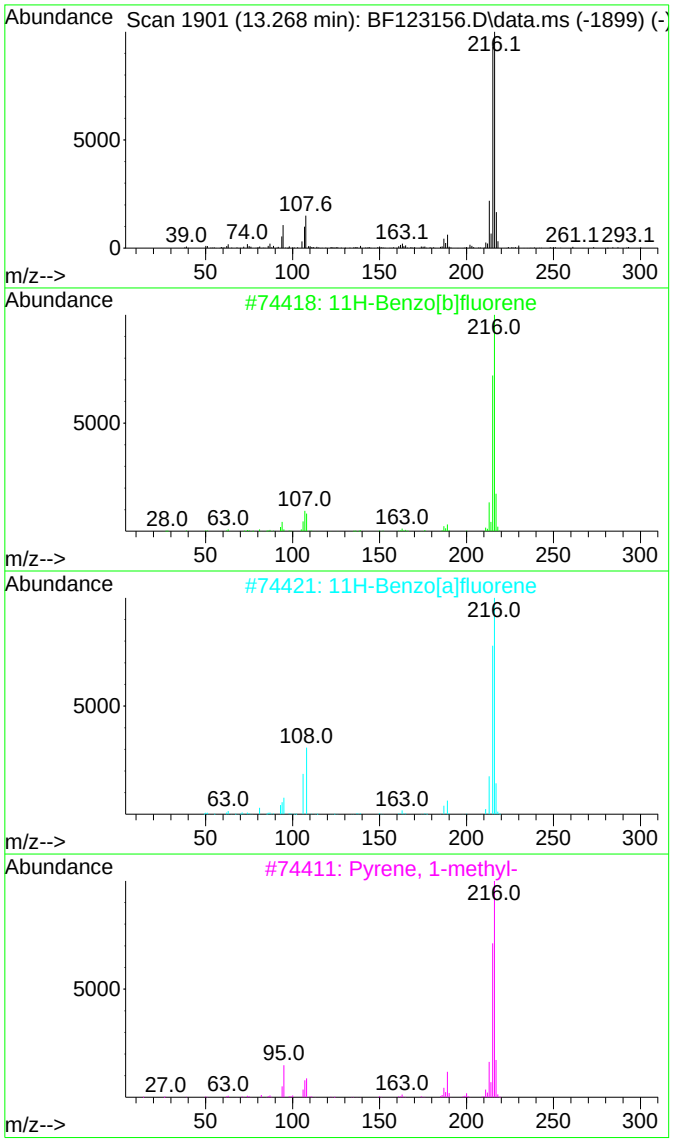
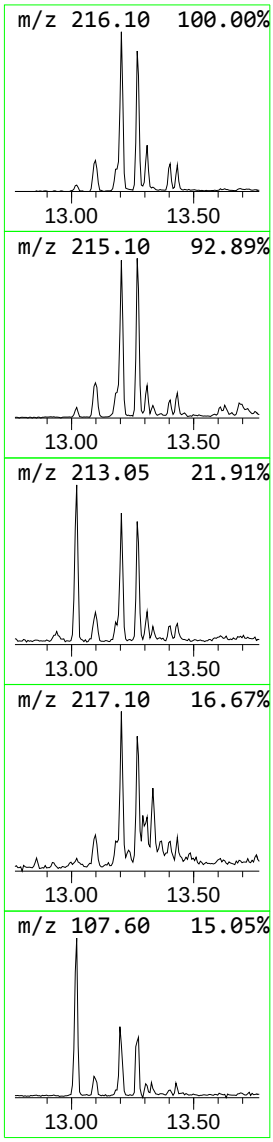
Peak Number 11 11H-Benzo[a]fluorene Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.269	2.25 ng	253244	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	93
2	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	87
3	Pyrene, 1-methyl-	216	C17H12	002381-21-7	80
4	7H-Benzo[c]fluorene	216	C17H12	000205-12-9	72
5	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	64





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Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
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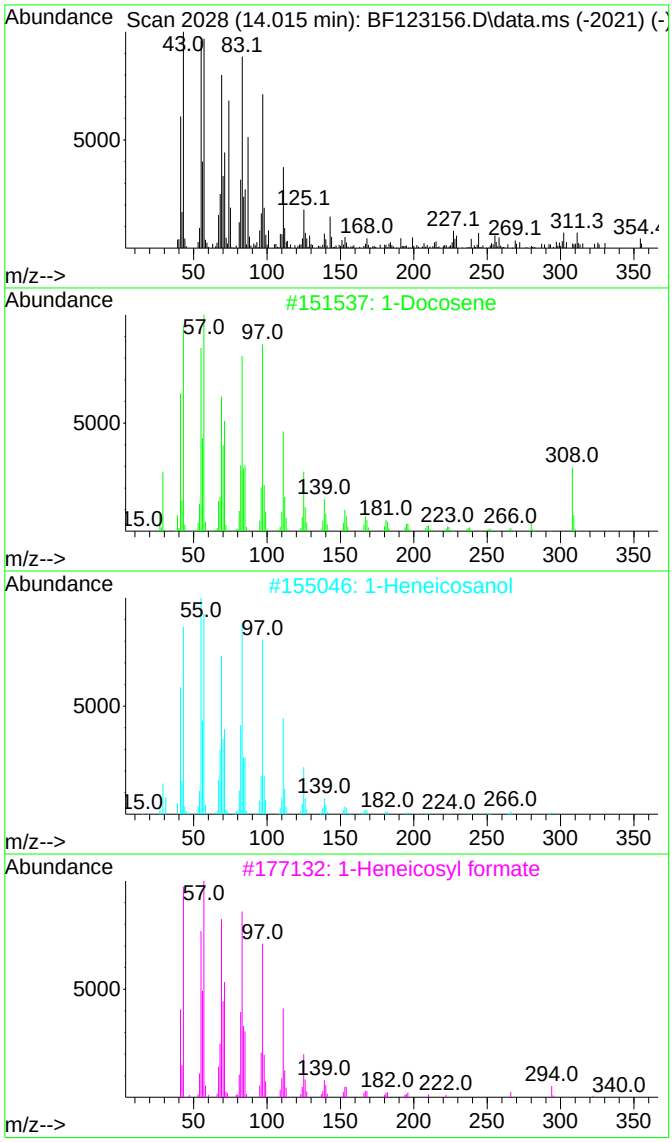
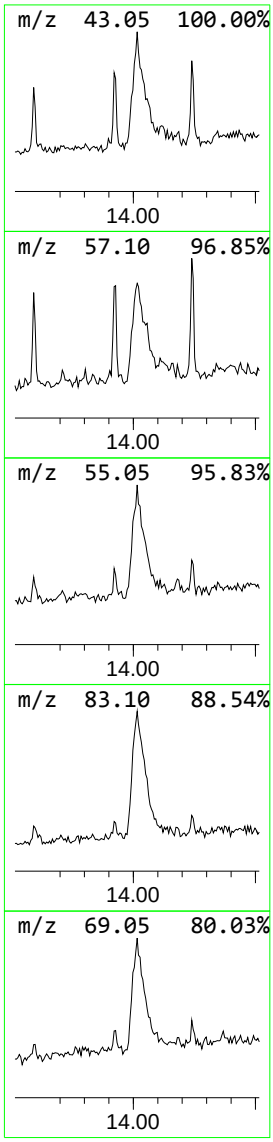
Peak Number 12 1-Docosene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.015	5.62 ng	632548	Chrysene-d12	14.080

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Docosene	308	C22H44	001599-67-3	95
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	1-Heneicosyl formate	340	C22H44O2	077899-03-7	93
4	n-Tetracosanol-1	354	C24H50O	000506-51-4	90
5	Trifluoroacetoxy hexadecane	338	C18H33F3O2	006222-03-3	86





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Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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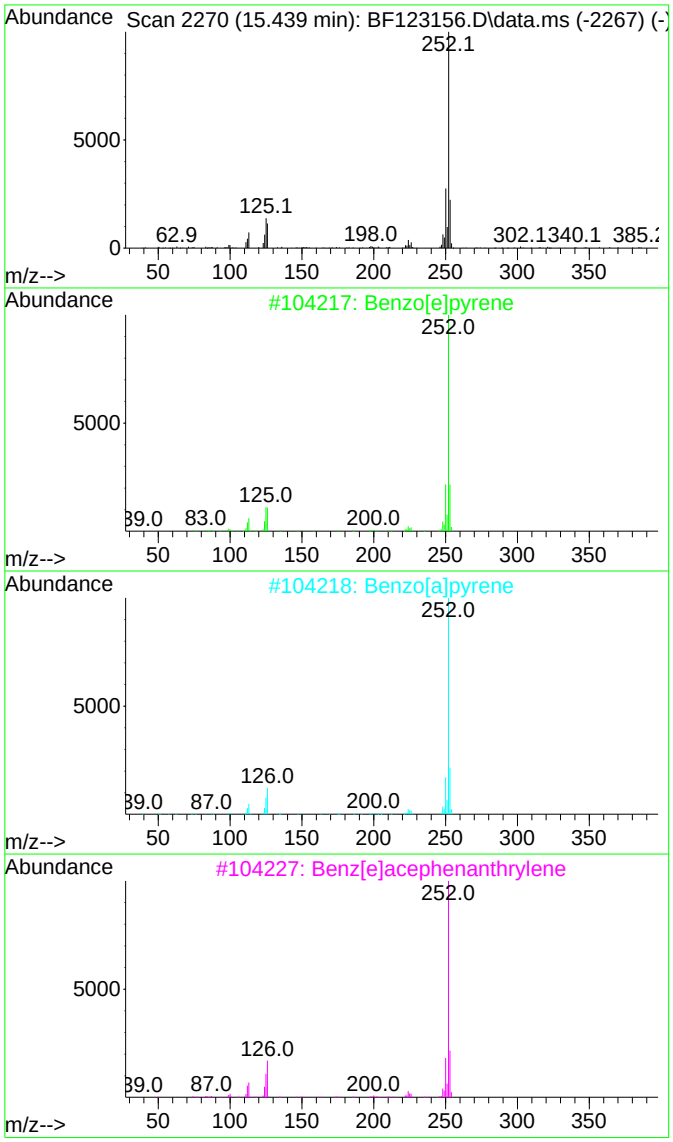
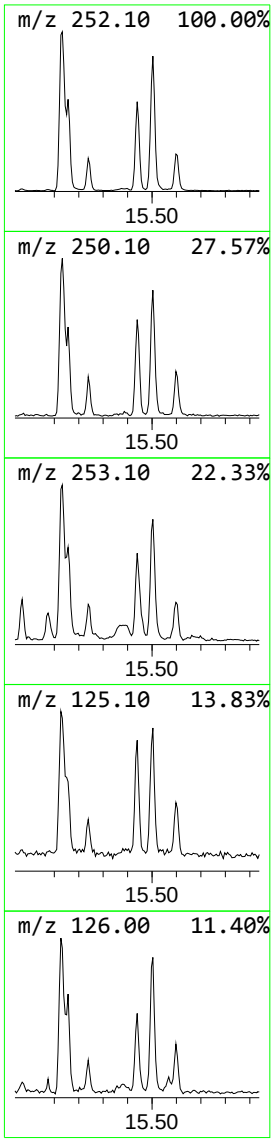
Peak Number 13 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.439	5.56 ng	339504	Perylene-d12	15.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	99
2	Benzo[a]pyrene	252	C20H12	000050-32-8	96
3	Benz[e]acephenanthrylene	252	C20H12	000205-99-2	95
4	Benzo[j]fluoranthene	252	C20H12	000205-82-3	95
5	Perylene	252	C20H12	000198-55-0	93





Tentatively Identified Compound (LSC) summary

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF020821\
Data File : BF123156.D
Acq On : 8 Feb 2021 21:59
Operator : JU/CG
Sample : M1332-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : TR-05-020521

Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF012721.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	2.181	11.3	ng	668892	1	6.892	1180090	20.0
5H-Indeno[1,2-b...	11.657	2.6	ng	239526	4	11.427	1845840	20.0
Hexadecanoic ac...	11.822	2.8	ng	261808	4	11.427	1845840	20.0
n-Hexadecanoic ...	11.957	6.1	ng	563838	4	11.427	1845840	20.0
4H-Cyclopenta[d...	12.045	3.3	ng	303070	4	11.427	1845840	20.0
9,12-Octadecadi...	12.510	9.5	ng	873376	4	11.427	1845840	20.0
11-Octadecenoic...	12.527	22.0	ng	2026850	4	11.427	1845840	20.0
Octadecanoic acid	12.745	10.4	ng	961098	4	11.427	1845840	20.0
11H-Benzo[b]flu...	13.204	2.6	ng	289935	5	14.080	2251850	20.0
11H-Benzo[a]flu...	13.269	2.3	ng	253244	5	14.080	2251850	20.0
1-Docosene	14.015	5.6	ng	632548	5	14.080	2251850	20.0
Benzo[e]pyrene	15.439	5.6	ng	339504	6	15.568	1222250	20.0