LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SB-21

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-BF110922.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

Signal : TIC: BF131138.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.152	5	11	17	rVB	453898	543967	38.29%	4.486%
2	5.104	509	513	522	rBV	193638	229394	16.15%	1.892%
3	5.504	577	581	585	rBV	1434529	1338291	94.20%	11.037%
4	6.516	748	753	756	rBV	1597812	1420749	100.00%	11.717%
5	6.886	813	816	820	rBV	507406	401636	28.27%	3.312%
6	7.451	908	912	915	rBV	1011710	883286	62.17%	7.285%
7	8.175	1031	1035	1038	rBV	514640	449414	31.63%	3.706%
8	9.245	1213	1217	1221	rBV	1309543	1191155	83.84%	9.824%
9	9.933	1330	1334	1337	rBV	433365	380864	26.81%	3.141%
10	10.722	1464	1468	1477	rBV	693586	562103	39.56%	4.636%
11	11.421	1584	1587	1590	rBV	427044	349731	24.62%	2.884%
12	11.927	1670	1673	1675	rBV	18925	17013	1.20%	0.140%
13	11.951	1675	1677	1681	rVB	25768	23051	1.62%	0.190%
14	12.004	1681	1686	1688	rBV2	12848	15191	1.07%	0.125%
15	12.033	1688	1691	1694	rVV2	37831	44284	3.12%	0.365%
16	12.063	1694	1696	1701	rVB	20685	17445	1.23%	0.144%
17	12.221	1715	1723	1726	rBV4	19235	28155	1.98%	0.232%
18	12.474	1763	1766	1769	rBV	62931	62972	4.43%	0.519%
19	12.510	1769	1772	1775	rVV2	29815	43046	3.03%	0.355%
20	12.563	1778	1781	1785	rBV3	26830	34456	2.43%	0.284%
21	12.639	1790	1794	1797	rBV	161543	130774	9.20%	1.079%
22	12.868	1827	1833	1839	rBV	222816	252602	17.78%	2.083%
23	12.921	1839	1842	1846	rVB3	30030	34773	2.45%	0.287%
24	13.010	1853	1857	1861	rBV	1203659	1064797	74.95%	8.781%
25	13.086	1867	1870	1877	rBV3	48261	71331	5.02%	0.588%
26	13.174	1882	1885	1886	rBV2	37786	37042	2.61%	0.305%
27	13.263	1898	1900	1903	rBV2	24344	22045	1.55%	0.182%
28	13.304	1903	1907	1910	rVV	67396	61095	4.30%	0.504%
29	13.398	1920	1923	1925	rBV	64395	52521	3.70%	0.433%
30	13.427	1925	1928	1931	rVB	58211	48428	3.41%	0.399%
31	13.710	1968	1976	1981	rBV2	42254	78063	5.49%	0.644%
32	13.821	1990	1995	1998	rBV4	36987	61640	4.34%	0.508%
33	13.851	1998	2000	2002	rVV	35265	28875	2.03%	0.238%
34	13.874	2002	2004	2007	rVV2	21629	23521	1.66%	0.194%
35	13.921	2009	2012	2015	rVB	24119	30507	2.15%	0.252%
36	14.074	2033	2038	2040	rBV2	545407	602695	42.42%	4.970%

LSC Area Percent Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

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-BF110922.M
Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

37	14.098	2040	2042	2049	rVB	182311	160219	11.28%	1.321%
38	14.168	2049	2054	2058	rBV3	59357	90345	6.36%	0.745%
39	14.227	2061	2064	2065	rBV3	16078	16485	1.16%	0.136%
40	14.327	2079	2081	2084	rVB4	16284	16855	1.19%	0.139%
41	14.421	2095	2097	2099	rBV	26162	24774	1.74%	0.204%
42	14.462	2099	2104	2107	rVV	75620	99040	6.97%	0.817%
43	14.492	2107	2109	2112	rVV	47907	40848	2.88%	0.337%
44	14.527	2113	2115	2118	rVV	33513	41693	2.93%	0.344%
45	14.586	2122	2125	2127	rBV2	36942	34949	2.46%	0.288%
46	15.127	2213	2217	2220	rBV	108623	150668	10.60%	1.243%
47	15.239	2233	2236	2240	rBV	25303	30778	2.17%	0.254%
48	15.439	2267	2270	2276	rVB	75854	92282	6.50%	0.761%
49	15.504	2277	2281	2287	rVB	100811	126870	8.93%	1.046%
50	15.568	2287	2292	2296	rBV	250599	335747	23.63%	2.769%
51	15.898	2345	2348	2350	rVB3	21162	18353	1.29%	0.151%
52	16.580	2461	2464	2465	rBV2	16454	16376	1.15%	0.135%
53	17.080	2543	2549	2557	rVB2	43226	96200	6.77%	0.793%
54	17.321	2587	2590	2596	rBV8	13502	28888	2.03%	0.238%
55	17.539	2622	2627	2634	rVB	30437	67260	4.73%	0.555%

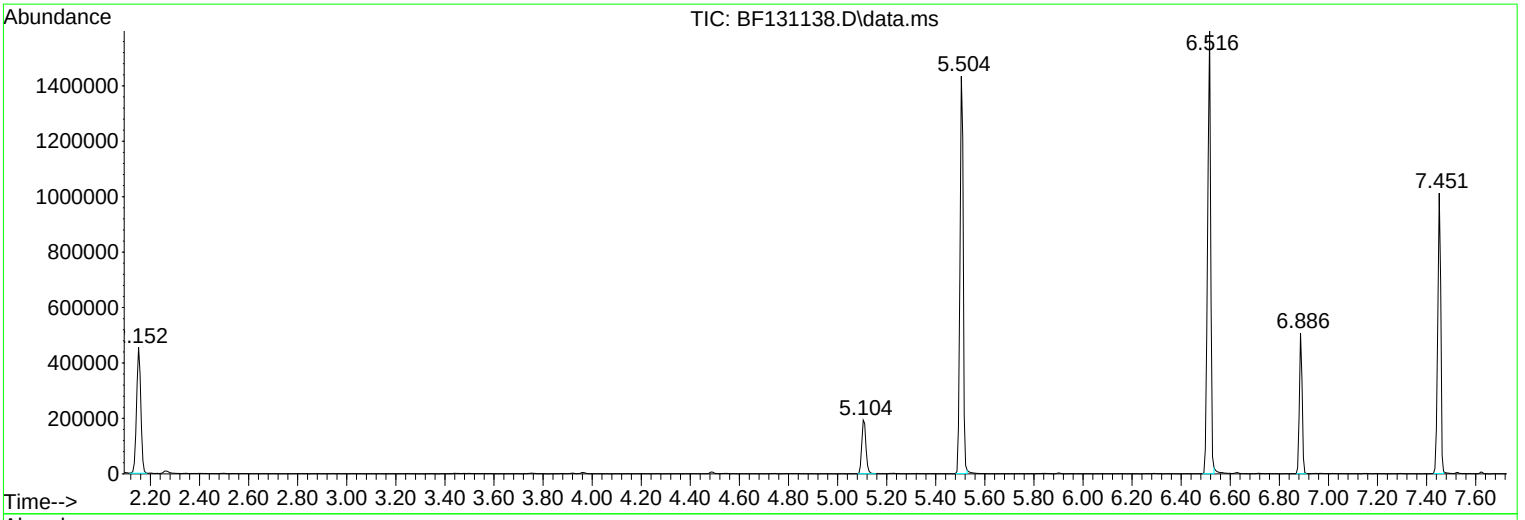
Sum of corrected areas: 12125542

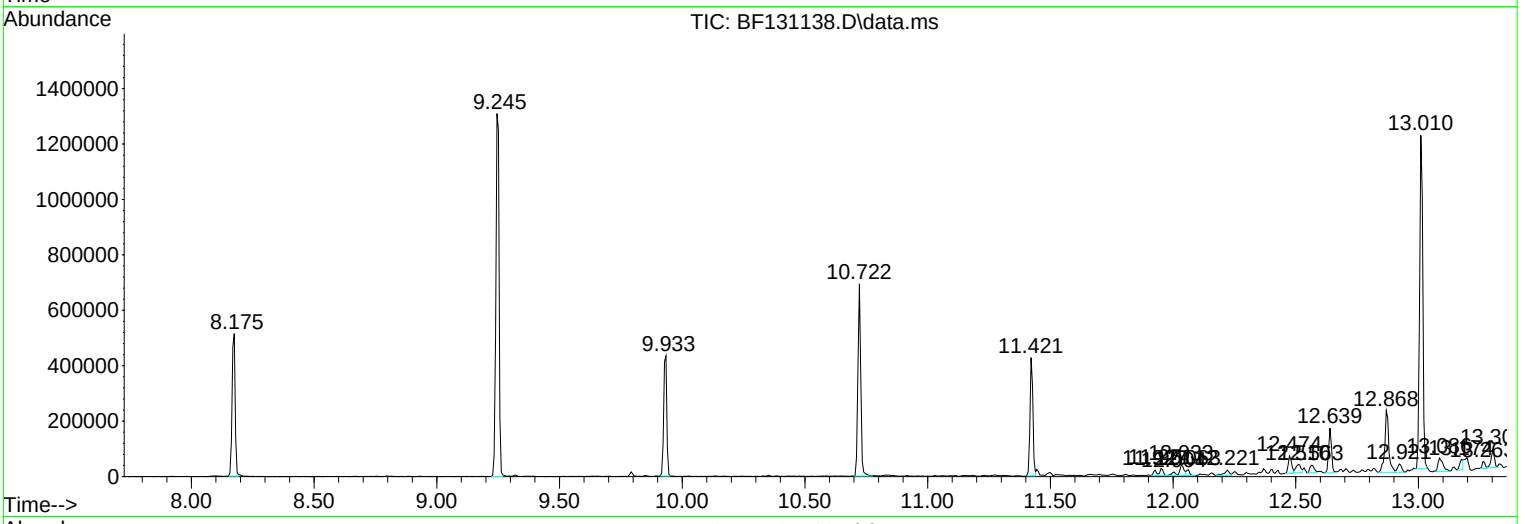
Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

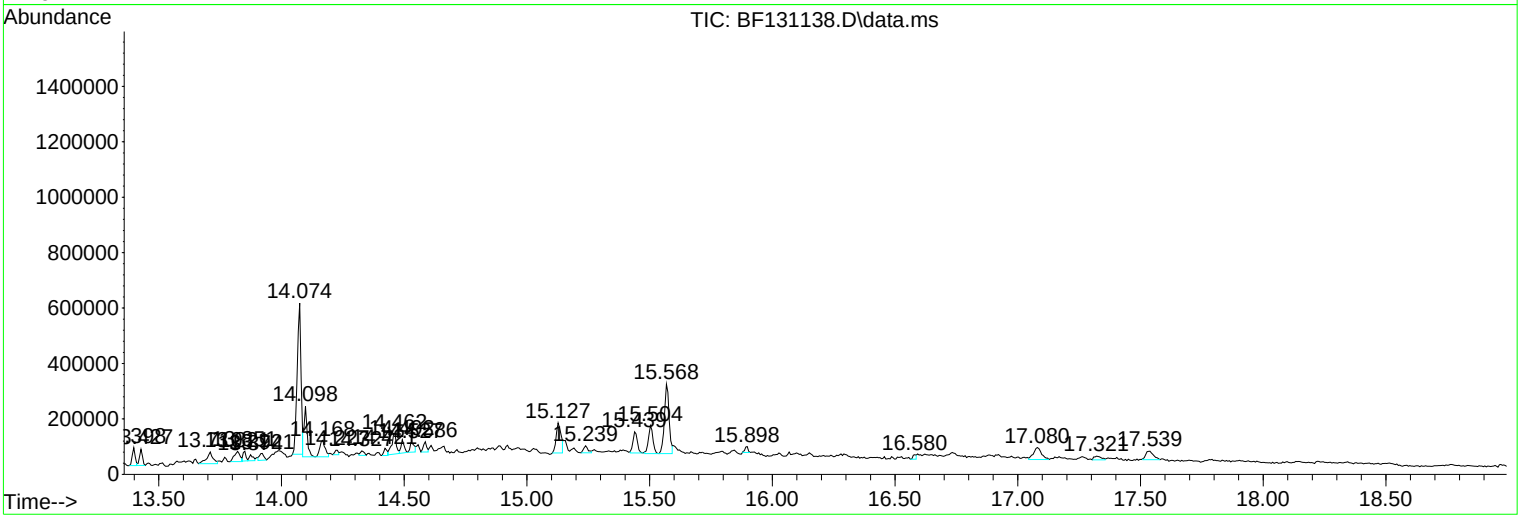
Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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







Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

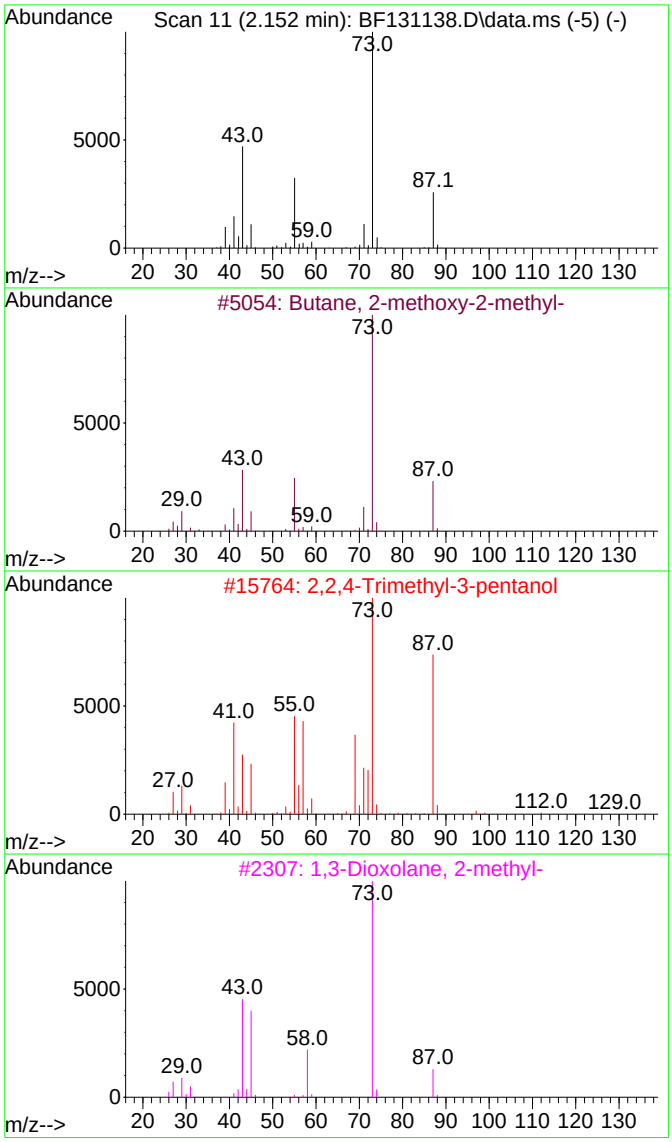
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

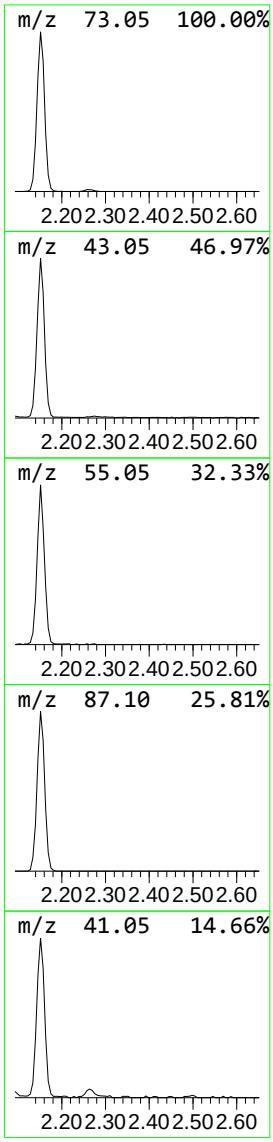
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
2.152	27.09 ng	543967	1,4-Dichlorobenzene-d4	6.886

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	40
3	1,3-Dioxolane, 2-methyl-	88	C4H8O2	000497-26-7	23
4	Octanal, 7-methoxy-3,7-dimethyl-	186	C11H22O2	003613-30-7	17
5	Pentane, 3-methoxy-	102	C6H14O	036839-67-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

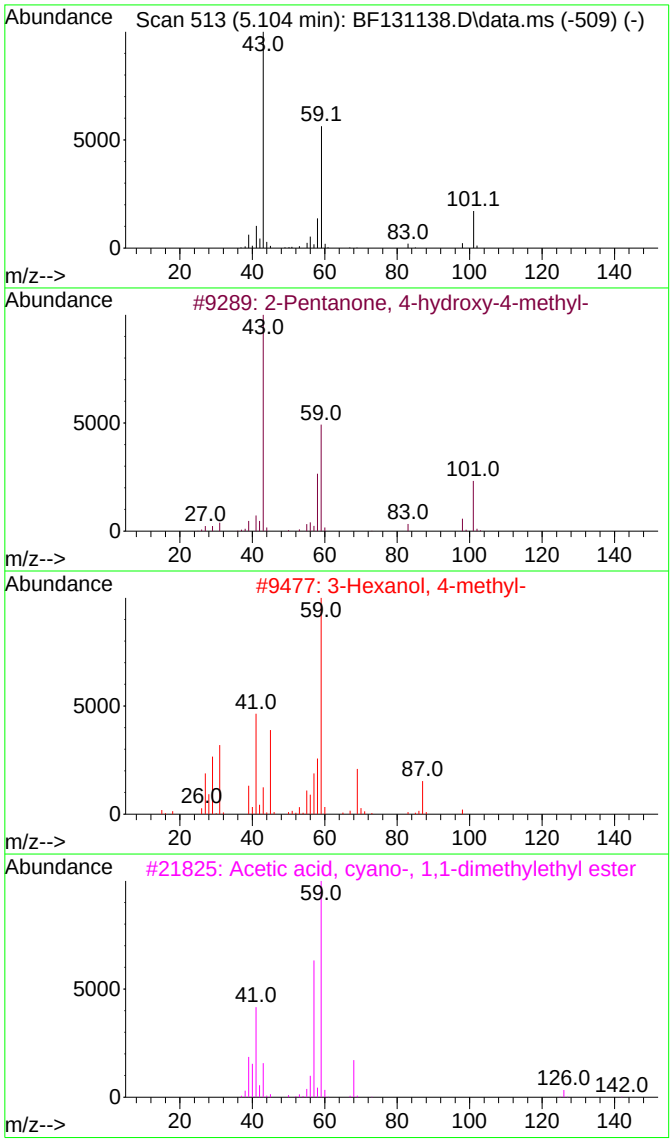
Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

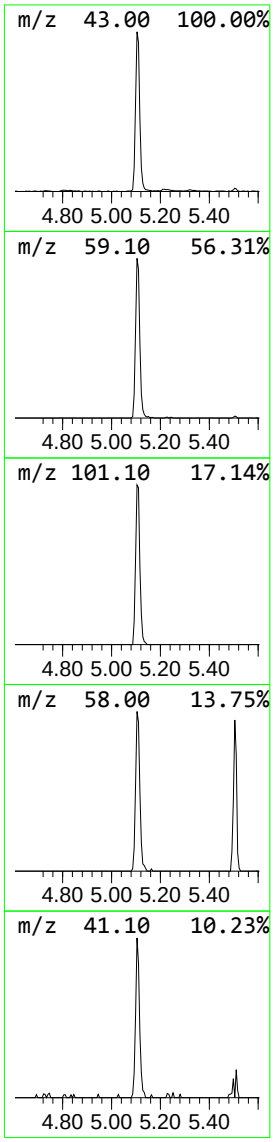
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

****
Peak Number 2 2-Pentanone, 4-hydroxy-4-me... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.104	11.42 ng	229394	1,4-Dichlorobenzene-d4	6.886

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Pentanone, 4-hydroxy-4-methyl-	116	C6H12O2	000123-42-2	38
2	3-Hexanol, 4-methyl-	116	C7H16O	000615-29-2	33
3	Acetic acid, cyano-, 1,1-dimethy...	141	C7H11NO2	001116-98-9	32
4	2,3-Butanedione, monooxime	101	C4H7NO2	000057-71-6	16
5	1-Propen-2-ol, acetate	100	C5H8O2	000108-22-5	12





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

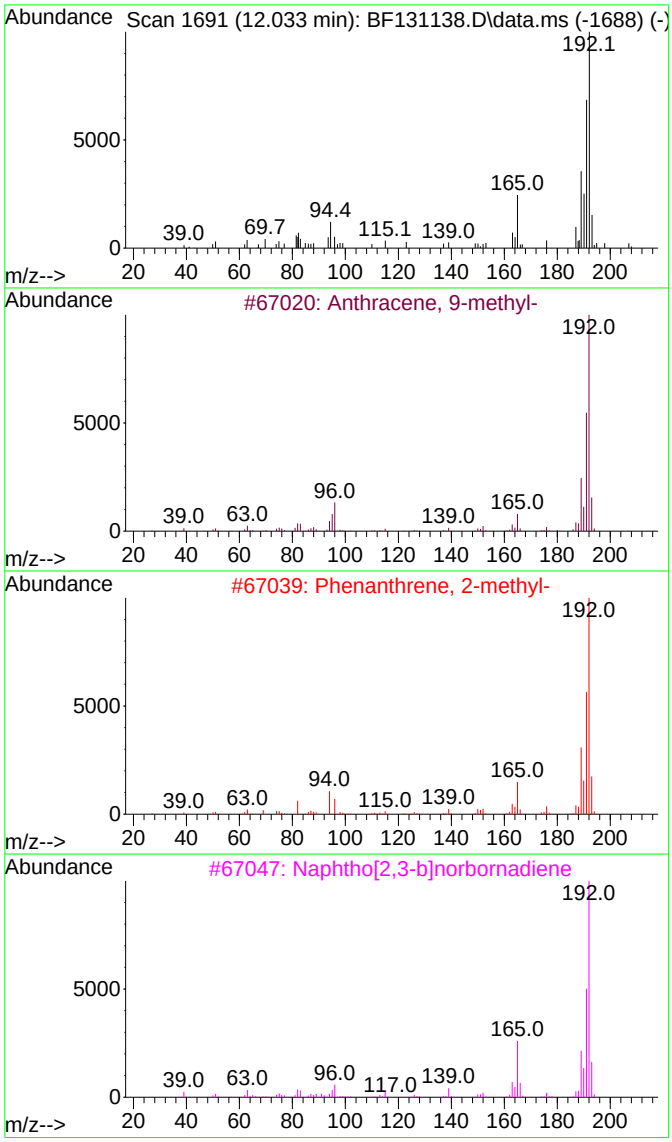
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

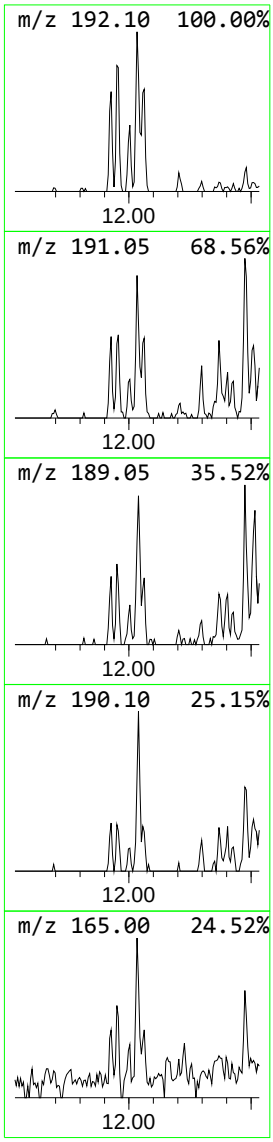
****

Peak Number 3 Anthracene, 9-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.033	2.53 ng	44284	Phenanthrene-d10	11.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Anthracene, 9-methyl-	192	C15H12	000779-02-2	81
2	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	76
3	Naphtho[2,3-b]norbornadiene	192	C15H12	107426-38-0	76
4	Anthracene, 1-methyl-	192	C15H12	000610-48-0	74
5	Phenanthrene, 4-methyl-	192	C15H12	000832-64-4	72





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

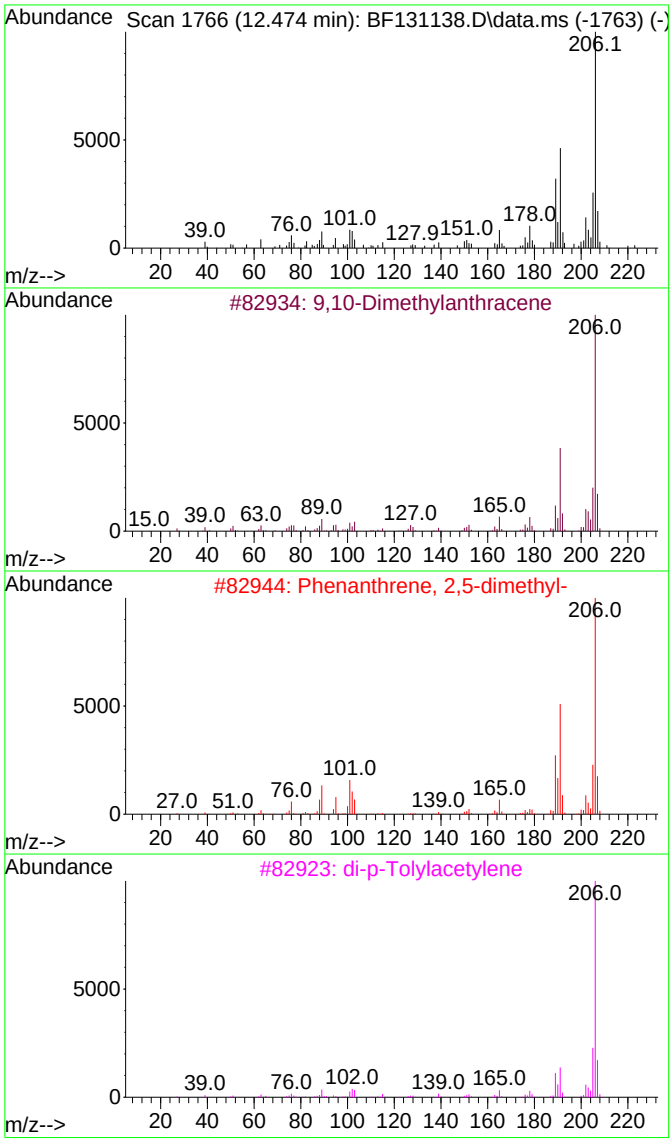
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

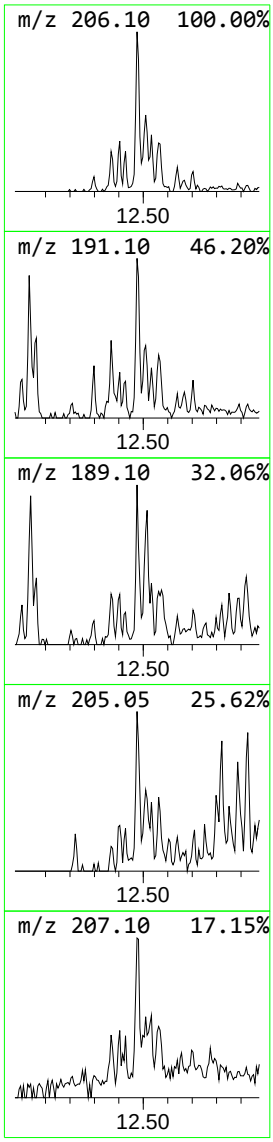
****

Peak Number 4 9,10-Dimethylanthracene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.474	3.60 ng	62972	Phenanthrene-d10	11.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,10-Dimethylanthracene	206	C16H14	000781-43-1	95
2	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	95
3	di-p-Tolylacetylene	206	C16H14	002789-88-0	93
4	Anthracene, 1,4-dimethyl-	206	C16H14	000781-92-0	91
5	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	91





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

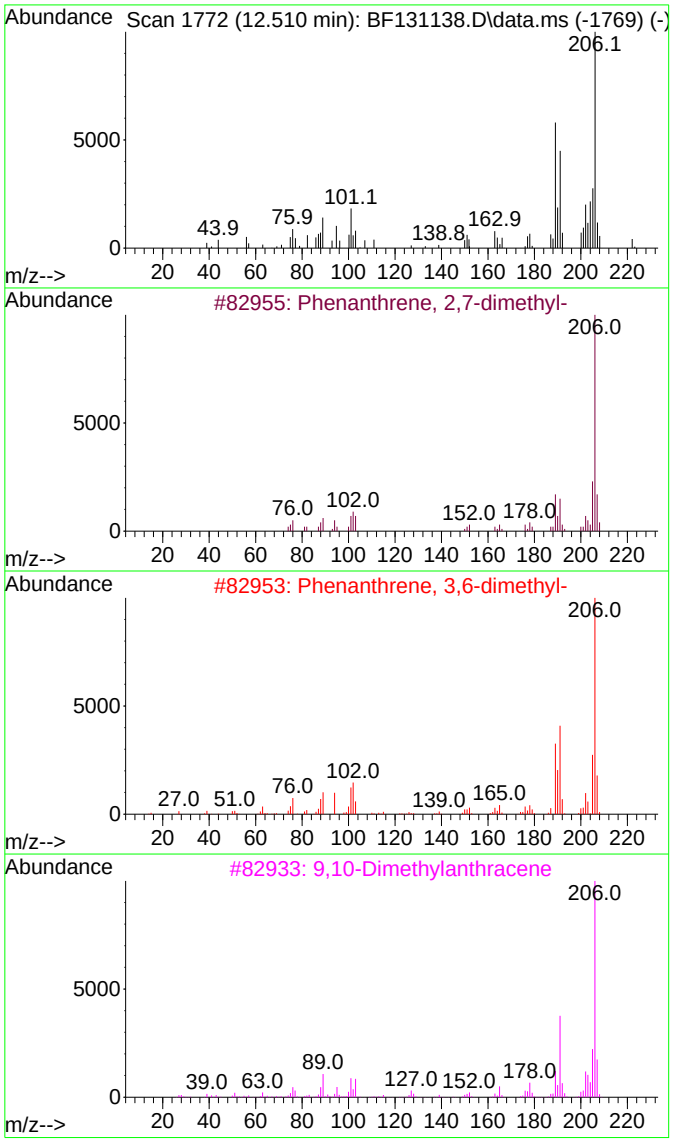
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

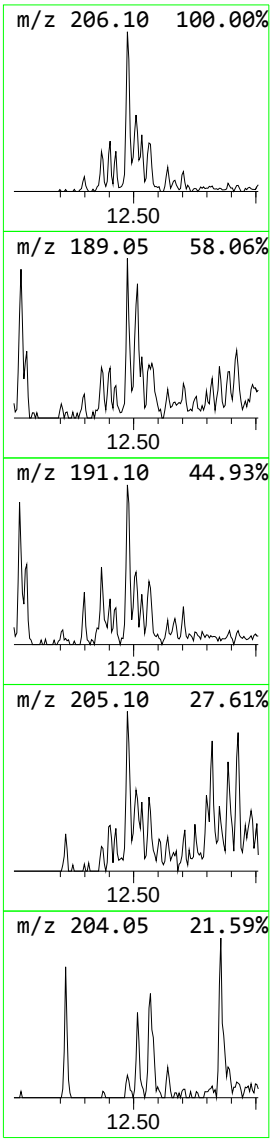
****

Peak Number 5 Phenanthrene, 2,7-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.510	2.46 ng	43046	Phenanthrene-d10	11.422

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	76
2	Phenanthrene, 3,6-dimethyl-	206	C16H14	001576-67-6	64
3	9,10-Dimethylanthracene	206	C16H14	000781-43-1	64
4	Naphthalene, 1,2-dihydro-4-phenyl-	206	C16H14	007469-40-1	60
5	Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	52





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Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

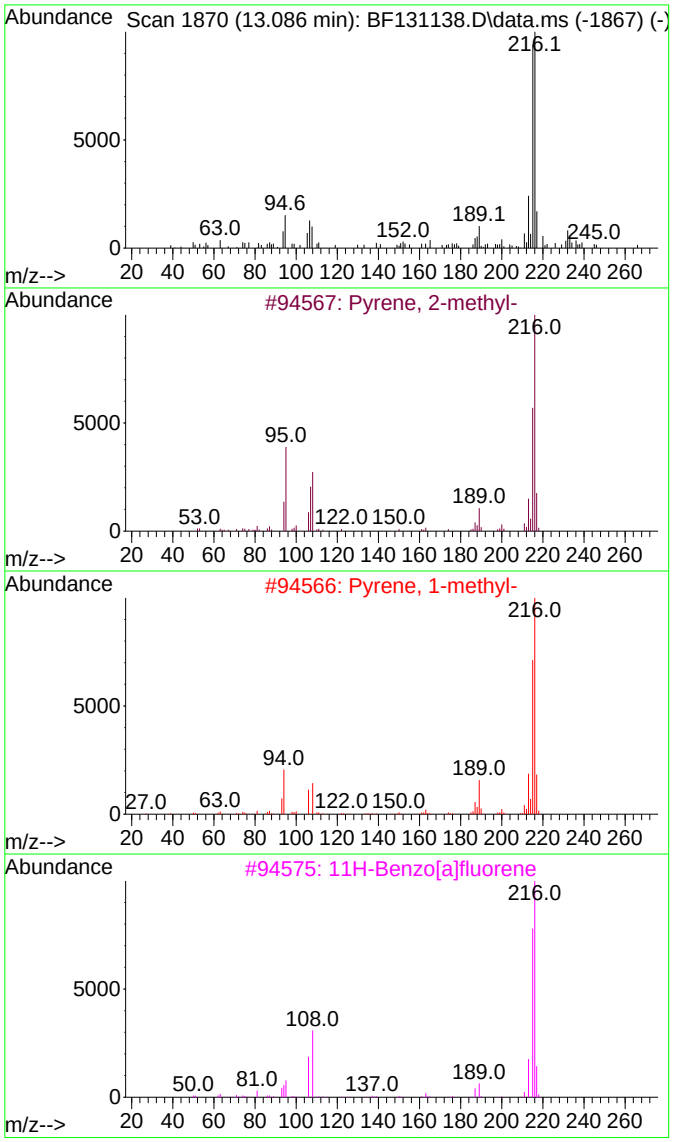
TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

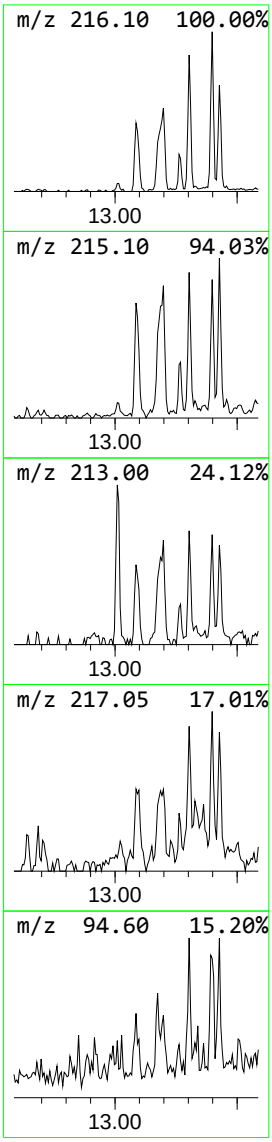
****

Peak Number 6 Pyrene, 2-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.086	2.37 ng	71331	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pyrene, 2-methyl-	216	C17H12	003442-78-2	91
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	89
3	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	83
4	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	81
5	11H-Benzo[b]fluorene	216	C17H12	000243-17-4	81





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.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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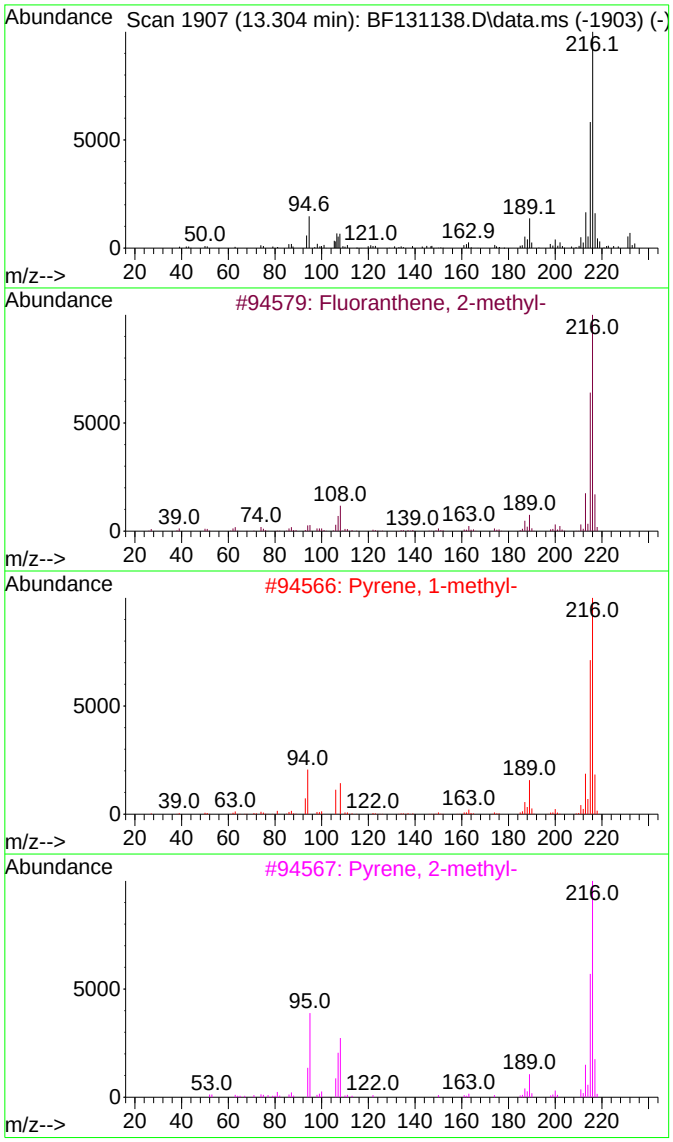
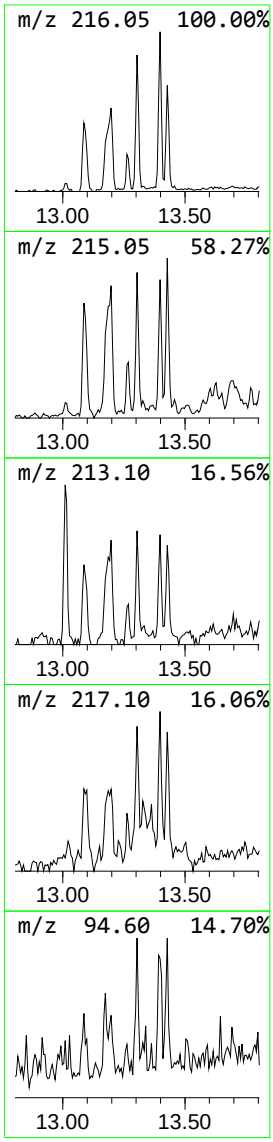
Peak Number 7 Fluoranthene, 2-methyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.304	2.03 ng	61095	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Fluoranthene, 2-methyl-	216	C17H12	033543-31-6	95
2	Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
3	Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
4	11H-Benzo[a]fluorene	216	C17H12	000238-84-6	90
5	Pyrene, 4-methyl-	216	C17H12	003353-12-6	89





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.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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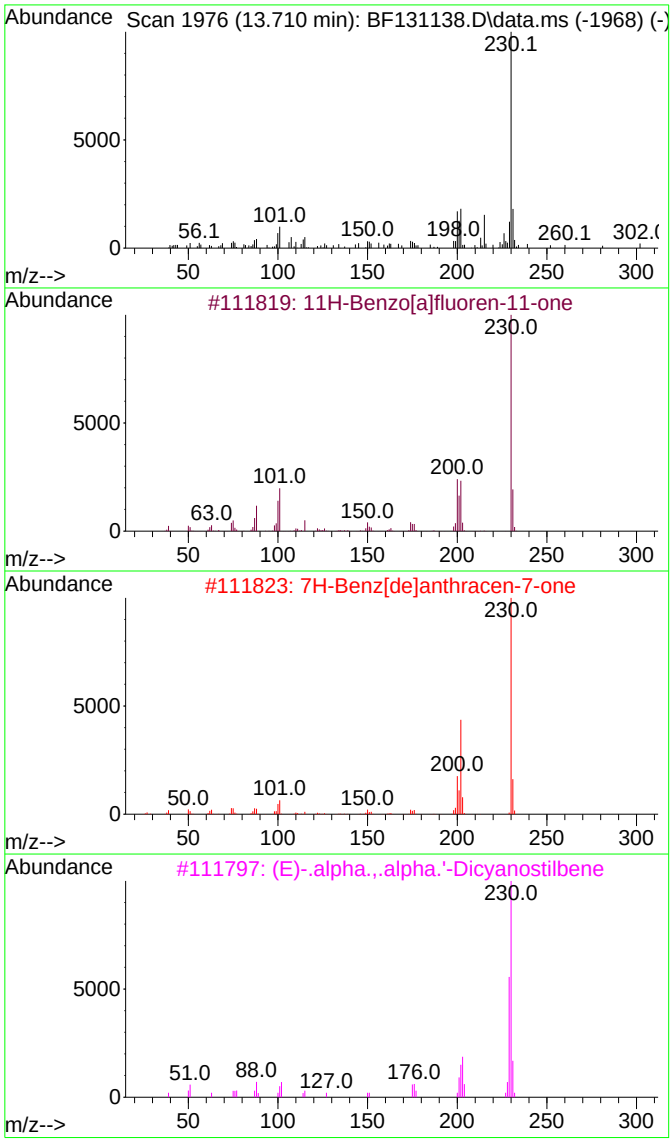
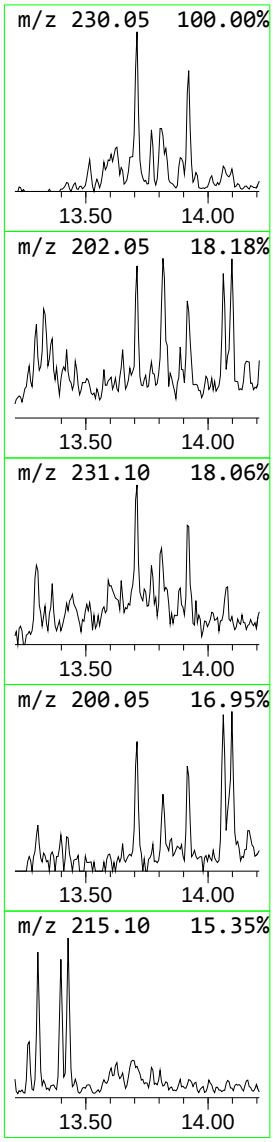
Peak Number 8 11H-Benzo[a]fluoren-11-one Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.710	2.59 ng	78063	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	11H-Benzo[a]fluoren-11-one	230	C17H10O	000479-79-8	94
2	7H-Benz[de]anthracen-7-one	230	C17H10O	000082-05-3	89
3	(E)-.alpha.,.alpha.'-Dicyanostil...	230	C16H10N2	002450-55-7	74
4	5-Amino-7-imino-4,8-diazatricycl...	230	C12H14N4O	1000388-35-6	68
5	4-Amino-1-imino-6-methyl-2H,6H,7...	230	C12H14N4O	1000411-06-8	64





Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF111022\
Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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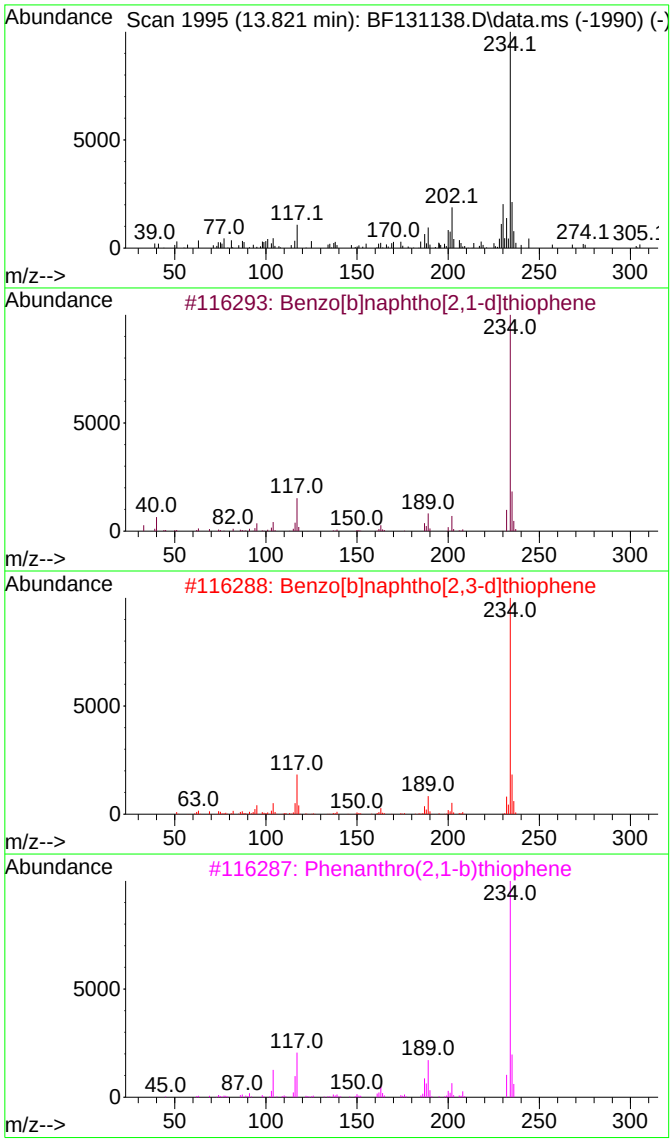
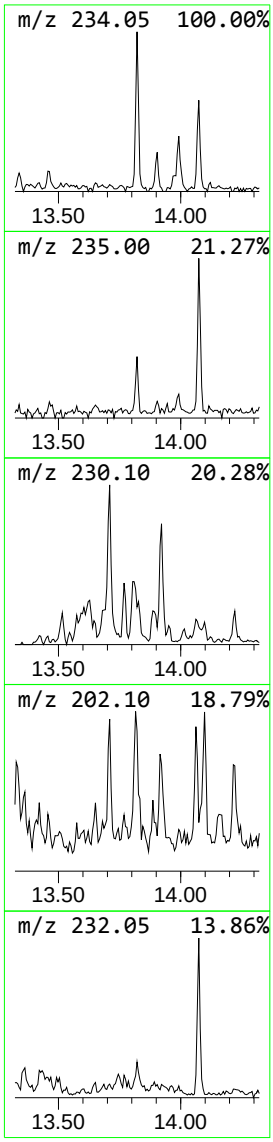
Peak Number 9 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.821	2.05 ng	61640	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	86
2	Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	62
3	Phenanthro(2,1-b)thiophene	234	C16H10S	000219-25-0	58
4	Benzo[b]naphtho[1,2-d]thiophene	234	C16H10S	000205-43-6	58
5	Quinoxalino[2,3-b]quinoxaline, 5...	234	C14H10N4	000531-46-4	50





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Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
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TIC Library : C:\Database\NIST20.L
TIC Integration Parameters: LSCINT.P

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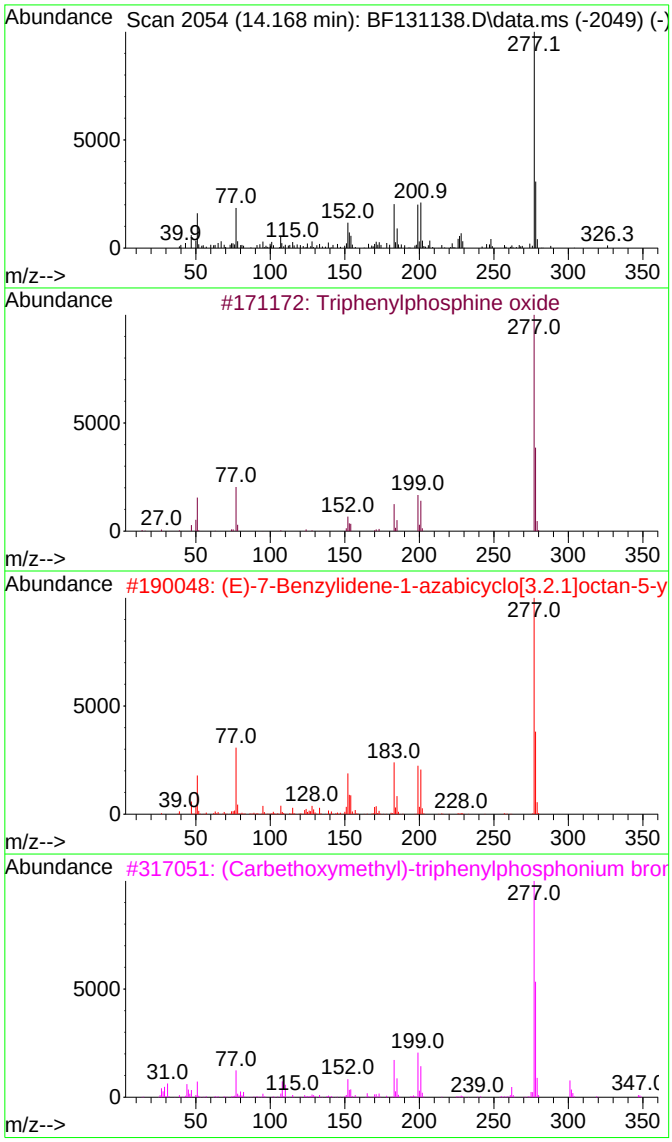
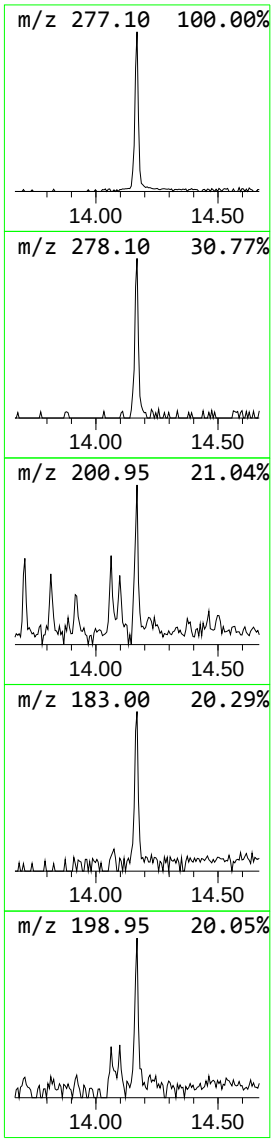
Peak Number 10 Triphenylphosphine oxide Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.168	3.00 ng	90345	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylphosphine oxide	278	C18H15OP	000791-28-6	97
2	(E)-7-Benzylidene-1-azabicyclo[3...	293	C15H19NO3S	1000483-83-4	58
3	(Carbethoxymethyl)-triphenylphos...	428	C22H22BrO2P	001530-45-6	50
4	5-(3,4-Dimethylphenylimino)-7ami...	277	C17H15N3O	1000254-01-4	38
5	5-Methyl-8-phenylfuro[3',2':5,6]...	277	C16H11N3O2	1000460-07-9	38





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Operator : CG\JU
Sample : N5494-19
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
SB-21

Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF110922.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

TIC Library : C:\Database\NIST20.L
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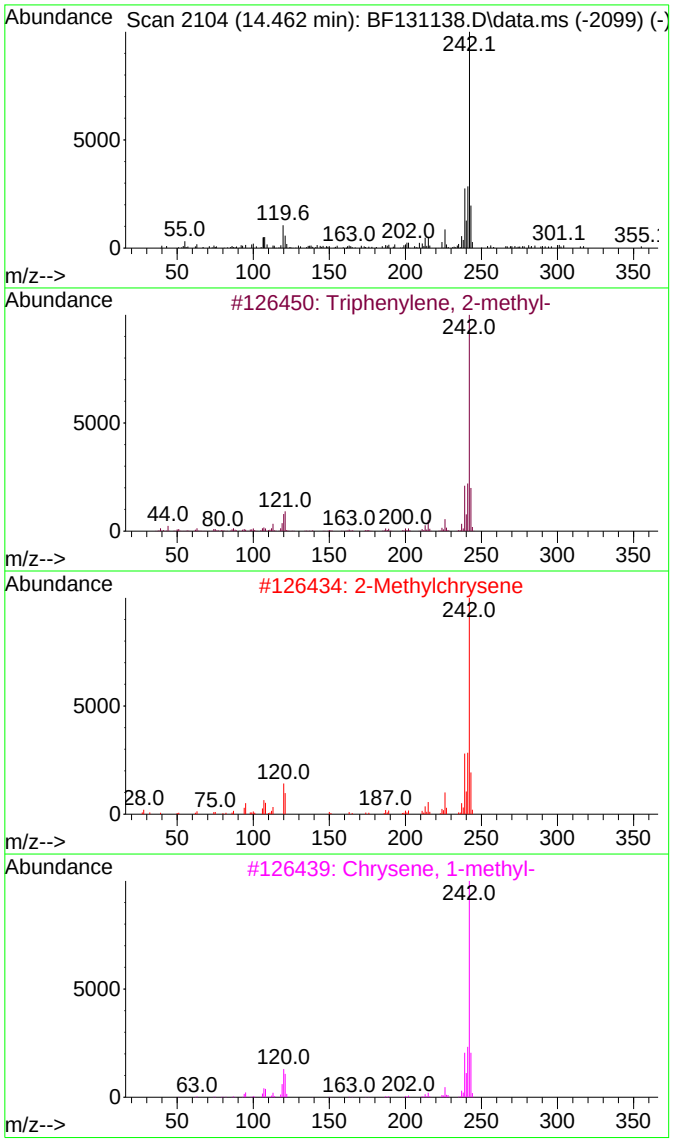
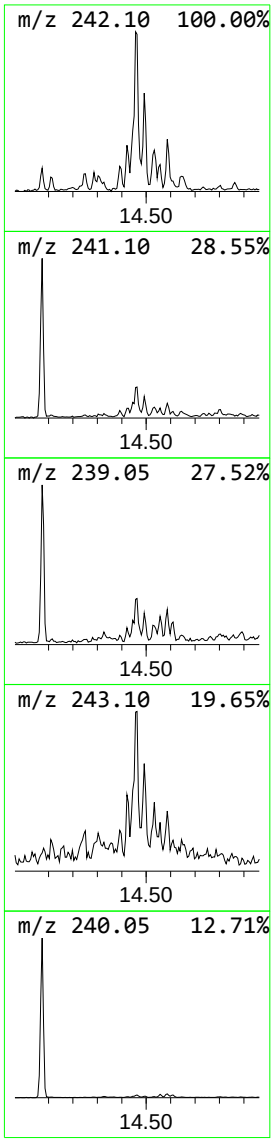
Peak Number 11 Triphenylene, 2-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	3.29 ng	99040	Chrysene-d12	14.074

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Triphenylene, 2-methyl-	242	C19H14	001705-84-6	93
2	2-Methylchrysene	242	C19H14	003351-32-4	93
3	Chrysene, 1-methyl-	242	C19H14	003351-28-8	91
4	Benz[a]anthracene, 2-methyl-	242	C19H14	002498-76-2	90
5	Benz[a]anthracene, 7-methyl-	242	C19H14	002541-69-7	89





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ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
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TIC Integration Parameters: LSCINT.P

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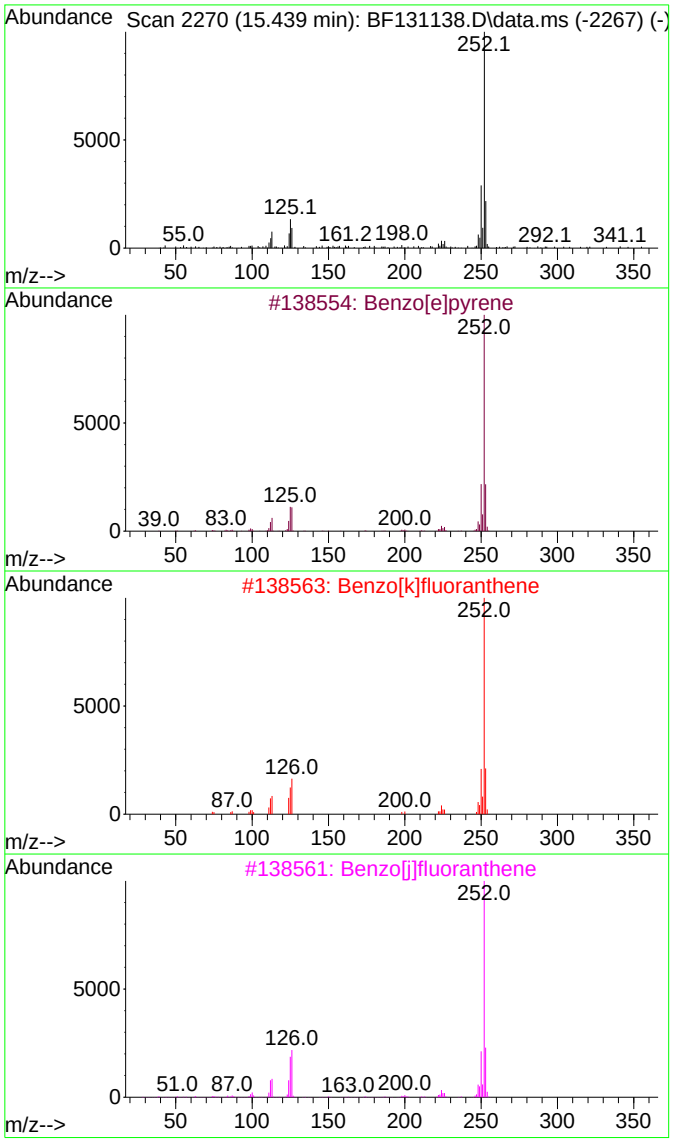
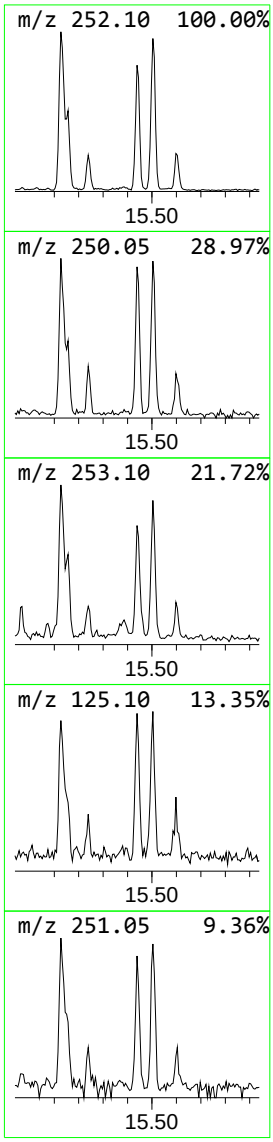
Peak Number 12 Benzo[e]pyrene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.439	5.50 ng	92282	Perylene-d12	15.568

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzo[e]pyrene	252	C20H12	000192-97-2	98
2	Benzo[k]fluoranthene	252	C20H12	000207-08-9	95
3	Benzo[j]fluoranthene	252	C20H12	000205-82-3	94
4	Perylene	252	C20H12	000198-55-0	91
5	Benzo[b]fluoranthene	252	C20H12	000205-99-2	90





Tentatively Identified Compound (LSC) summary

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Data File : BF131138.D
Acq On : 10 Nov 2022 23:13
Operator : CG\JU
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Misc :
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Instrument :
BNA_F
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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

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

					--Internal Standard---			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
Butane, 2-metho...	2.152	27.1	ng	543967	1	6.886	401636	20.0
2-Pentanone, 4-...	5.104	11.4	ng	229394	1	6.886	401636	20.0
Anthracene, 9-m...	12.033	2.5	ng	44284	4	11.422	349731	20.0
9,10-Dimethylan...	12.474	3.6	ng	62972	4	11.422	349731	20.0
Phenanthrene, 2...	12.510	2.5	ng	43046	4	11.422	349731	20.0
Pyrene, 2-methyl-	13.086	2.4	ng	71331	5	14.074	602695	20.0
Fluoranthene, 2...	13.304	2.0	ng	61095	5	14.074	602695	20.0
11H-Benzo[a]flu...	13.710	2.6	ng	78063	5	14.074	602695	20.0
Benzo[b]naphtho...	13.821	2.0	ng	61640	5	14.074	602695	20.0
Triphenylphosph...	14.168	3.0	ng	90345	5	14.074	602695	20.0
Triphenylene, 2...	14.463	3.3	ng	99040	5	14.074	602695	20.0
Benzo[e]pyrene	15.439	5.5	ng	92282	6	15.568	335747	20.0