

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
 Data File : BF089919.D
 Acq On : 26 Aug 2016 20:36
 Operator : UM/SJ
 Sample : PB93109BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93109BS

Quant Time: Aug 26 23:17:29 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 18:51:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.65	152	316943	20.00	ng	0.00
21) Naphthalene-d8	7.94	136	1275329	20.00	ng	0.00
38) Acenaphthene-d10	9.68	164	584434	20.00	ng	-0.01
63) Phenanthrene-d10	11.15	188	1053165	20.00	ng	0.00
75) Chrysene-d12	13.77	240	727066	20.00	ng	-0.01
86) Perylene-d12	15.12	264	641600	20.00	ng	-0.01

System Monitoring Compounds						
5) 2-Fluorophenol	5.23	112	1866852	101.01	ng	0.01
7) Phenol-d6	6.30	99	2107690	92.90	ng	0.00
23) Nitrobenzene-d5	7.22	82	1346283	65.58	ng	0.00
41) 2,4,6-Tribromophenol	10.47	330	481196	86.64	ng	-0.01
44) 2-Fluorobiphenyl	9.02	172	2213919	66.58	ng	0.00
78) Terphenyl-d14	12.74	244	2013156	62.64	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.19	88	252462	27.60	ng	96
3) Pyridine	2.82	79	729447	30.41	ng	93
4) n-Nitrosodimethylamine	2.79	42	308357	34.53	ng	# 83
6) Aniline	6.31	93	651940	21.42	ng	# 9
8) 2-Chlorophenol	6.43	128	736639	33.33	ng	89
9) Benzaldehyde	6.18	77	136235	9.22	ng	97
10) Phenol	6.31	94	776185	29.19	ng	99
11) bis(2-Chloroethyl)ether	6.40	93	701933	34.04	ng	90
12) 1,3-Dichlorobenzene	6.59	146	776872m	33.62	ng	
13) 1,4-Dichlorobenzene	6.67	146	760628m	32.03	ng	
14) 1,2-Dichlorobenzene	6.82	146	753772	34.00	ng	96
15) Benzyl Alcohol	6.80	79	481060	31.97	ng	95
16) 2,2'-oxybis(1-Chloropropan	6.95	45	822873	31.18	ng	90
17) 2-Methylphenol	6.92	107	611000	35.30	ng	94
18) Hexachloroethane	7.16	117	278326	32.86	ng	94
19) n-Nitroso-di-n-propylamine	7.08	70	499412	34.42	ng	# 83
20) 3+4-Methylphenols	7.07	107	655430	30.88	ng	# 74
22) Acetophenone	7.07	105	873062	29.81	ng	# 88
24) Nitrobenzene	7.23	77	639755	27.78	ng	99
25) Isophorone	7.48	82	1331591	32.16	ng	# 93
26) 2-Nitrophenol	7.55	139	354679	30.87	ng	# 83
27) 2,4-Dimethylphenol	7.61	122	665415	32.13	ng	95
28) bis(2-Chloroethoxy)methane	7.70	93	909411	35.50	ng	97
29) 2,4-Dichlorophenol	7.79	162	563170	32.41	ng	90
30) 1,2,4-Trichlorobenzene	7.88	180	633292	32.95	ng	# 95
31) Naphthalene	7.95	128	1944104	32.62	ng	99
32) Benzoic acid	7.72	122	330408	20.47	ng	96
33) 4-Chloroaniline	8.01	127	208876	8.08	ng	# 95
34) Hexachlorobutadiene	8.08	225	310513	30.85	ng	99
35) Caprolactam	8.38	113	169135m	31.96	ng	
36) 4-Chloro-3-methylphenol	8.50	107	617239	32.80	ng	99
37) 2-Methylnaphthalene	8.65	142	1401204	36.34	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.82	216	450924	23.83	ng	97
40) Hexachlorocyclopentadiene	8.81	237	606499	92.31	ng	98

Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\
 Data File : BF089919.D
 Acq On : 26 Aug 2016 20:36
 Operator : UM/SJ
 Sample : PB93109BS
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB93109BS

Quant Time: Aug 26 23:17:29 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 18:51:25 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.92	196	393566	33.06	nq	98
43) 2,4,5-Trichlorophenol	8.97	196	417471	35.25	nq	99
45) 1,1'-Biphenyl	9.12	154	1397140	30.89	nq	93
46) 2-Chloronaphthalene	9.13	162	1144530	32.01	nq	98
47) 2-Nitroaniline	9.23	65	394011	38.64	nq	# 74
48) Acenaphthylene	9.54	152	1764988	32.85	nq	100
49) Dimethylphthalate	9.43	163	1479519	35.95	nq	# 98
50) 2,6-Dinitrotoluene	9.47	165	295254	31.08	nq	88
51) Acenaphthene	9.71	154	1278206	35.95	nq	99
52) 3-Nitroaniline	9.63	138	187168	17.78	nq	96
53) 2,4-Dinitrophenol	9.75	184	334747	65.18	nq	90
54) Dibenzofuran	9.88	168	1594632	35.71	nq	97
55) 4-Nitrophenol	9.82	139	624192	72.65	nq	# 71
56) 2,4-Dinitrotoluene	9.87	165	383885	31.14	nq	# 64
57) Fluorene	10.23	166	1115303	31.34	nq	98
58) 2,3,4,6-Tetrachlorophenol	10.01	232	349002m	40.49	nq	
59) Diethylphthalate	10.12	149	1487062	35.57	nq	97
60) 4-Chlorophenyl-phenylether	10.23	204	527977	32.52	nq	99
61) 4-Nitroaniline	10.25	138	322644	32.70	nq	98
62) Azobenzene	10.39	77	1535548	39.07	nq	92
64) 4,6-Dinitro-2-methylphenol	10.28	198	236678	39.11	nq	# 60
65) n-Nitrosodiphenylamine	10.34	169	1162041	35.09	nq	99
66) 4-Bromophenyl-phenylether	10.71	248	354698	32.12	nq	# 67
67) Hexachlorobenzene	10.78	284	433783	34.44	nq	98
68) Atrazine	10.88	200	362246	35.72	nq	97
69) Pentachlorophenol	10.97	266	491242	68.80	nq	98
70) Phenanthrene	11.18	178	1805576	33.94	nq	99
71) Anthracene	11.23	178	2002055	37.03	nq	99
72) Carbazole	11.38	167	1786952	36.85	nq	99
73) Di-n-butylphthalate	11.74	149	2197925	36.22	nq	99
74) Fluoranthene	12.35	202	1757741	34.74	nq	94
76) Benzidine	12.48	184	659999	28.05	nq	97
77) Pyrene	12.58	202	1827921	30.87	nq	98
79) Butylbenzylphthalate	13.22	149	972364	36.36	nq	98
80) Benzo(a)anthracene	13.76	228	1517960	36.46	nq	99
81) 3,3'-Dichlorobenzidine	13.74	252	262439	19.23	nq	# 95
82) Chrysene	13.81	228	1453044	40.70	nq	99
83) Bis(2-ethylhexyl)phthalate	13.78	149	1204240	32.69	nq	# 98
84) Di-n-octyl phthalate	14.39	149	2011509	41.10	nq	# 92
85) Indeno(1,2,3-cd)pyrene	16.37	276	1254321	27.54	nq	99
87) Benzo(b)fluoranthene	14.77	252	1549460m	43.90	nq	
88) Benzo(k)fluoranthene	14.79	252	969919m	30.29	nq	
89) Benzo(a)pyrene	15.07	252	1171869	35.08	nq	# 96
90) Dibenzo(a,h)anthracene	16.39	278	1053139	29.04	nq	97
91) Benzo(g,h,i)perylene	16.74	276	1081322	28.32	nq	# 94

(#) = qualifier out of range (m) = manual integration (+) = signals summed

```
Data Path : Z:\HPCHEM1\BNA F\DATA\BF082616\  
Data File : BF089919.D  
Acq On    : 26 Aug 2016   20:36  
Operator  : UM/SJ  
Sample    : PB93109BS  
Misc      :  
ALS Vial  : 8      Sample Multiplier: 1
```

Instrument :
BNA_F
ClientSampleId :
PB93109BS

Quant Time: Aug 26 23:17:29 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081916.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 26 18:51:25 2016
Response via : Initial Calibration

