

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050115\
 Data File : BF078896.D
 Acq On : 1 May 2015 16:54
 Operator : TP/IZ
 Sample : PB83053BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB83053BS

Quant Time: May 02 00:34:00 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 02 00:19:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	252136m	20.00	ng	0.18
21) Naphthalene-d8	0.00	136	0	0.00	ng	-8.94
38) Acenaphthene-d10	0.00	164	0	0.00	ng	-11.12
63) Phenanthrene-d10	0.00	188	0	0.00	ng	-12.96
75) Chrysene-d12	16.26	240	538	20.00	ng	-0.02
86) Perylene-d12	18.07	264	911	20.00	ng	-0.06
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	553609	37.80	ng	0.00
7) Phenol-d6	6.90	99	723417	37.36	ng	0.00
23) Nitrobenzene-d5	8.04	82	392092	0.00	ng	0.00
41) 2,4,6-Tribromophenol	12.09	330	218673	0.00	ng	0.00
44) 2-Fluorobiphenyl	10.28	172	863456	0.00	ng	0.00
78) Terphenyl-d14	14.96	244	1052475	46836.04	ng	0.00
Target Compounds						
2) 1,4-Dioxane	2.34	88	57191	8.98	ng	Qvalue # 32
3) Pyridine	3.11	79	179151	9.61	ng	96
4) n-Nitrosodimethylamine	3.02	42	74577	9.34	ng	98
6) Aniline	6.95	93	167859	6.14	ng	99
8) 2-Chlorophenol	7.08	128	183074	11.02	ng	98
9) Benzaldehyde	6.81	77	77341	5.93	ng	98
10) Phenol	6.92	94	248583	11.07	ng	93
11) bis(2-Chloroethyl)ether	7.04	93	166096	10.09	ng	98
12) 1,3-Dichlorobenzene	7.28	146	188421	10.09	ng	98
13) 1,4-Dichlorobenzene	7.38	146	191762	9.98	ng	98
14) 1,2-Dichlorobenzene	7.56	146	188050	10.51	ng	99
15) Benzyl Alcohol	7.53	79	153851	10.70	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.71	45	266283	10.57	ng	99
17) 2-Methylphenol	7.67	107	144966	10.84	ng	96
18) Hexachloroethane	7.99	117	65796	9.94	ng	93
19) n-Nitroso-di-n-propylamine	7.87	70	134808	10.31	ng	# 83
20) 3+4-Methylphenols	7.86	107	198292	11.13	ng	89
76) Benzidine	14.64	184	240023	16553.75	ng	98
77) Pyrene	14.75	202	672316	21686.18	ng	99
79) Butylbenzylphthalate	15.58	149	311866	23012.87	ng	# 88
80) Benzo(a)anthracene	16.25	228	622603	21132.72	ng	100
81) 3,3'-Dichlorobenzidine	16.23	252	169598	16161.18	ng	# 97
82) Chrysene	16.30	228	590169	20827.51	ng	100
83) Bis(2-ethylhexyl)phthalate	16.31	149	460251	22422.51	ng	99
84) Di-n-octyl phthalate	17.13	149	791998	21908.17	ng	# 100
85) Indeno(1,2,3-cd)pyrene	19.58	276	754036	20885.22	ng	# 100
87) Benzo(b)fluoranthene	17.60	252	695159	13941.16	ng	99
88) Benzo(k)fluoranthene	17.63	252	535724	11097.97	ng	98
89) Benzo(a)pyrene	18.00	252	607932	13240.78	ng	# 96
90) Dibenzo(a,h)anthracene	19.61	278	609688	12494.50	ng	97
91) Benzo(g,h,i)perylene	20.03	276	625235	12784.17	ng	99

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

Data Path : Z:\HPCHEM1\BNA F\DATA\BF050115\
Data File : BF078896.D
Acq On : 1 May 2015 16:54
Operator : TP/IZ
Sample : PB83053BS
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_F
ClientSampleId :
PB83053BS

Quant Time: May 02 00:34:00 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 02 00:19:57 2015
Response via : Initial Calibration

