

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121516\
 Data File : BF091667.D
 Acq On : 16 Dec 2016 10:40
 Operator : UM/SJ
 Sample : H6006-22DL 5X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 WC-2-0-5DL

Quant Time: Dec 16 16:56:47 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 16 00:35:41 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.78	152	51448	20.00	ng	0.00
21) Naphthalene-d8	8.06	136	198173	20.00	ng	-0.01
38) Acenaphthene-d10	9.82	164	103191	20.00	ng	0.00
63) Phenanthrene-d10	11.30	188	110168	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	130196	20.00	ng	0.00
86) Perylene-d12	15.35	264	110444	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	103163	33.78	ng	-0.01
7) Phenol-d6	6.42	99	226674	59.54	ng	-0.01
23) Nitrobenzene-d5	7.34	82	132591	37.34	ng	-0.01
41) 2,4,6-Tribromophenol	10.61	330	51102	59.62	ng	0.00
44) 2-Fluorobiphenyl	9.15	172	290395	48.63	ng	0.00
78) Terphenyl-d14	12.89	244	215383	37.54	ng	-0.01

Target Compounds				Qvalue		
10) Phenol	6.43	94	16438	3.69	ng	87
31) Naphthalene	8.09	128	29048	3.15	ng	99
32) Benzoic acid	7.73	122	2300	5.89	ng	# 7
48) Acenaphthylene	9.69	152	40041	4.54	ng	98
49) Dimethylphthalate	9.54	163	50861	7.78	ng	97
51) Acenaphthene	9.85	154	31662	5.17	ng	97
54) Dibenzofuran	10.03	168	36384	4.80	ng	# 88
55) 4-Nitrophenol	10.03	139	12993	10.37	ng	# 1
65) n-Nitrosodiphenylamine	10.48	169	7387	2.25	ng	96
70) Phenanthrene	11.32	178	458506	84.04	ng	99
71) Anthracene	11.38	178	110590	18.79	ng	98
72) Carbazole	11.53	167	62819	12.18	ng	99
74) Fluoranthene	12.51	202	636384	111.40	ng	97
77) Pyrene	12.74	202	536494	51.94	ng	98
79) Butylbenzylphthalate	13.37	149	10085	2.50	ng	97
80) Benzo(a)anthracene	13.93	228	445064	58.86	ng	98
82) Chrysene	13.93	228	445064	56.26	ng	99
83) Bis(2-ethylhexyl)phthalate	13.93	149	16998	3.29	ng	# 96
85) Indeno(1,2,3-cd)pyrene	16.69	276	153358	21.54	ng	95
87) Benzo(b)fluoranthene	14.95	252	605505	91.97	ng	99
88) Benzo(k)fluoranthene	14.95	252	605505	104.60	ng	100
89) Benzo(a)pyrene	15.29	252	269455	45.43	ng	# 96
90) Dibenzo(a,h)anthracene	16.71	278	42010	7.91	ng	# 89
91) Benzo(g,h,i)perylene	17.09	276	134922	24.95	ng	# 91

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

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