

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101254.D
 Acq On : 8 Dec 2017 23:34
 Operator : SJ/JU
 Sample : I6615-01
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PB-2-OW(0-10)

Quant Time: Dec 09 02:31:47 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF113017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 08 18:17:31 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.83	152	44085	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	163510	20.00	ng	0.00
38) Acenaphthene-d10	9.87	164	69816	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	121346	20.00	ng	0.00
75) Chrysene-d12	14.00	240	98440	20.00	ng	0.00
86) Perylene-d12	15.48	264	81531	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	299242	112.17	ng	0.02
7) Phenol-d6	6.47	99	356500	110.06	ng	0.01
23) Nitrobenzene-d5	7.39	82	200998	73.33	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	82087	97.88	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	422770	82.85	ng	0.00
78) Terphenyl-d14	12.95	244	363419	80.75	ng	0.00
Target Compounds						
10) Phenol	6.48	94	13986	3.883	ng	Qvalue # 63
49) Dimethylphthalate	9.57	163	19508	3.465	ng	99
70) Phenanthrene	11.38	178	45253	6.772	ng	97
74) Fluoranthene	12.57	202	74661	10.079	ng	94
77) Pyrene	12.80	202	73058	10.755	ng	99
80) Benzo(a)anthracene	14.00	228	34506	5.552	ng	97
82) Chrysene	14.03	228	31912	5.528	ng	96
83) Bis(2-ethylhexyl)phthalate	13.97	149	8718	2.042	ng	# 89
85) Indeno(1,2,3-cd)pyrene	16.96	276	15185	2.959	ng	# 90
87) Benzo(b)fluoranthene	15.05	252	32608m	6.677	ng	
88) Benzo(k)fluoranthene	15.07	252	10156m	2.111	ng	
89) Benzo(a)pyrene	15.42	252	25682	5.785	ng	# 90
91) Benzo(a,h,i)perylene	17.42	276	15667	4.247	ng	# 90

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

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