

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101245.D
 Acq On : 8 Dec 2017 19:31
 Operator : SJ/JU
 Sample : I6684-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-01-120617

Quant Time: Dec 08 22:14:34 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	6.82	152	48482	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	190677	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	82511	20.00	ng	0.00
63) Phenanthrene-d10	11.35	188	139101	20.00	ng	0.00
75) Chrysene-d12	14.00	240	101952	20.00	ng	0.00
86) Perylene-d12	15.47	264	107389	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	303975	103.61	ng	0.00
7) Phenol-d6	6.46	99	363259	101.98	ng	0.00
23) Nitrobenzene-d5	7.39	82	219491	68.67	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	87477	88.25	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	377836	62.65	ng	0.00
78) Terphenyl-d14	12.94	244	274126	58.81	ng	0.00
Target Compounds						
10) Phenol	6.47	94	14157	3.574	ng	99
49) Dimethylphthalate	9.57	163	44748	6.725	ng	97
70) Phenanthrene	11.38	178	118337	15.448	ng	95
71) Anthracene	11.43	178	32359	4.174	ng	97
74) Fluoranthene	12.57	202	274587	32.338	ng	95
77) Pyrene	12.80	202	243735	34.646	ng	98
80) Benzo(a)anthracene	13.99	228	148460	23.063	ng	100
82) Chrysene	14.03	228	113064	18.913	ng	96
85) Indeno(1,2,3-cd)pyrene	16.94	276	52489	9.876	ng	# 91
87) Benzo(b)fluoranthene	15.04	252	151072m	23.486	ng	
88) Benzo(k)fluoranthene	15.07	252	41991m	6.626	ng	
89) Benzo(a)pyrene	15.41	252	94134m	16.097	ng	
90) Dibenzo(a,h)anthracene	16.95	278	17661m	3.426	ng	
91) Benzo(a,h,i)perylene	17.40	276	47887	9.856	ng	# 91

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

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