

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101263.D
 Acq On : 9 Dec 2017 5:14
 Operator : SJ/JU
 Sample : I6791-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AREA-2-1-SAMPLE

Quant Time: Dec 11 00:24:52 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	41523	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	151297	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	63768	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	111530	20.00	ng	0.00
75) Chrysene-d12	14.01	240	84924	20.00	ng	0.00
86) Perylene-d12	15.48	264	68626	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.45	112	135964	54.11	ng	0.00
7) Phenol-d6	6.46	99	327379	107.31	ng	0.00
23) Nitrobenzene-d5	7.39	82	189547	74.73	ng	0.00
41) 2,4,6-Tribromophenol	10.66	330	18839	24.59	ng	0.00
44) 2-Fluorobiphenyl	9.19	172	410896	88.16	ng	0.00
78) Terphenyl-d14	12.94	244	346968	89.36	ng	0.00
Target Compounds						
10) Phenol	6.47	94	13041	3.844	ng	82
49) Dimethylphthalate	9.57	163	40157	7.808	ng	# 99
70) Phenanthrene	11.38	178	13086	2.131	ng	97
74) Fluoranthene	12.57	202	32027	4.704	ng	94
77) Pyrene	12.80	202	32715	5.583	ng	97
80) Benzo(a)anthracene	14.00	228	17941	3.346	ng	97
82) Chrysene	14.03	228	16498	3.313	ng	93
85) Indeno(1,2,3-cd)pyrene	16.96	276	10632	2.402	ng	# 91
87) Benzo(b)fluoranthene	15.05	252	20819m	5.065	ng	
89) Benzo(a)pyrene	15.42	252	15342	4.105	ng	# 81
91) Benzo(g,h,i)perylene	17.41	276	11929	3.842	ng	# 90

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

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