

Data Path : Z:\HPCHEM1\BNA F\DATA\BF120817\
 Data File : BF101248.D
 Acq On : 8 Dec 2017 20:52
 Operator : SJ/JU
 Sample : I6677-03 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-120717-B

Quant Time: Dec 08 22:53:52 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	6.82	152	44838	20.00	ng	0.00
21) Naphthalene-d8	8.11	136	173694	20.00	ng	0.00
38) Acenaphthene-d10	9.86	164	72825	20.00	ng	0.00
63) Phenanthrene-d10	11.36	188	119173	20.00	ng	0.00
75) Chrysene-d12	14.00	240	97491	20.00	ng	0.00
86) Perylene-d12	15.47	264	80935	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.44	112	56892	20.97	ng	0.00
7) Phenol-d6	6.45	99	68862	20.90	ng	-0.01
23) Nitrobenzene-d5	7.39	82	39667	13.62	ng	-0.01
41) 2,4,6-Tribromophenol	10.66	330	14546	16.63	ng	0.00
44) 2-Fluorobiphenyl	9.18	172	79983	15.03	ng	0.00
78) Terphenyl-d14	12.93	244	59106	13.26	ng	0.00
Target Compounds						
70) Phenanthrene	11.38	178	60280	9.185	ng	96
71) Anthracene	11.43	178	15043	2.265	ng	98
74) Fluoranthene	12.57	202	84484	11.613	ng	93
77) Pyrene	12.80	202	71954	10.696	ng	98
80) Benzo(a)anthracene	13.99	228	30316m	4.925	ng	
82) Chrysene	14.03	228	29855	5.222	ng	96
85) Indeno(1,2,3-cd)pyrene	16.94	276	11249	2.213	ng	91
87) Benzo(b)fluoranthene	15.04	252	30925m	6.379	ng	
89) Benzo(a)pyrene	15.41	252	21453	4.868	ng	95
91) Benzo(g,h,i)perylene	17.40	276	11112	3.035	ng	# 86

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

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