

Data Path : Z:\HPCHEM1\BNA F\DATA\BF011215\
 Data File : BF076826.D
 Acq On : 13 Jan 2015 3:01
 Operator : IZ / TP
 Sample : G1062-03
 Misc : S
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW-013

Quant Time: Jan 13 10:23:24 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF010915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 12 10:43:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	34222	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	153557	20.00	ng	-0.01
38) Acenaphthene-d10	15.12	164	81918	20.00	ng	-0.01
63) Phenanthrene-d10	17.85	188	128742	20.00	ng	0.00
75) Chrysene-d12	21.34	240	121311	20.00	ng	-0.01
86) Perylene-d12	22.99	264	111832	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.17	112	190383	92.83	ng	-0.01
7) Phenol-d6	7.48	99	243463	94.30	ng	-0.02
23) Nitrobenzene-d5	9.37	82	145960	52.82	ng	-0.01
41) 2,4,6-Tribromophenol	16.77	330	53893	78.82	ng	-0.01
44) 2-Fluorobiphenyl	13.64	172	287420	51.64	ng	-0.01
78) Terphenyl-d14	20.07	244	255430	46.23	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	14.70	163	30691	5.16	ng	97
70) Phenanthrene	17.89	178	138542	17.27	ng	99
71) Anthracene	17.97	178	31112	3.99	ng	99
74) Fluoranthene	19.52	202	154320	18.78	ng	96
77) Pyrene	19.81	202	161569	18.51	ng	96
80) Benzo(a)anthracene	21.33	228	65247	8.44	ng	98
82) Chrysene	21.37	228	50850	7.27	ng	96
85) Indeno(1,2,3-cd)pyrene	24.09	276	31790m	4.59	ng	
87) Benzo(b)fluoranthene	22.57	252	69144m	8.93	ng	
88) Benzo(k)fluoranthene	22.61	252	17076m	2.52	ng	
89) Benzo(a)pyrene	22.93	252	49306	7.41	ng	94
91) Benzo(a,h,i)perylene	24.40	276	30714	4.96	ng	96

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

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