

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF121515\
 Data File : BF083814.D
 Acq On : 15 Dec 2015 12:53
 Operator : UM/IZ
 Sample : G4634-01
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 K203-0-2

Quant Time: Dec 16 00:57:35 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF121315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 14 01:30:12 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	129857	20.00	ng	-0.02
21) Naphthalene-d8	8.18	136	528310	20.00	ng	-0.03
38) Acenaphthene-d10	9.94	164	259502	20.00	ng	-0.02
63) Phenanthrene-d10	11.43	188	423836	20.00	ng	-0.02
75) Chrysene-d12	14.05	240	321664	20.00	ng	-0.02
86) Perylene-d12	15.49	264	294353	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	874331	108.62	ng	-0.01
7) Phenol-d6	6.53	99	950643	93.26	ng	-0.01
23) Nitrobenzene-d5	7.46	82	601277	63.67	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	255753	96.72	ng	-0.02
44) 2-Fluorobiphenyl	9.26	172	1069088	54.50	ng	-0.02
78) Terphenyl-d14	13.00	244	631853	42.25	ng	-0.02
Target Compounds						Qvalue
10) Phenol	6.54	94	31081	2.65	ng	# 72
48) Acenaphthylene	9.80	152	90405	3.17	ng	98
49) Dimethylphthalate	9.65	163	278165	13.75	ng	99
70) Phenanthrene	11.45	178	115341	5.12	ng	98
74) Fluoranthene	12.63	202	228213	9.81	ng	95
77) Pyrene	12.85	202	253987	9.99	ng	99
80) Benzo(a)anthracene	14.04	228	122704	6.71	ng	# 86
82) Chrysene	14.08	228	136046m	7.63	ng	
85) Indeno(1,2,3-cd)pyrene	16.91	276	114319	6.52	ng	# 100
87) Benzo(b)fluoranthene	15.08	252	205381m	10.94	ng	
88) Benzo(k)fluoranthene	15.09	252	53322m	3.13	ng	
89) Benzo(a)pyrene	15.43	252	136080	8.12	ng	# 93
91) Benzo(g,h,i)perylene	17.35	276	130032	7.59	ng	97

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

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