

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
 Data File : BF078839.D
 Acq On : 30 Apr 2015 5:38
 Operator : TP/IZ
 Sample : G2021-04
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-17-3-5

Quant Time: Apr 30 07:24:11 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 30 01:56:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	79848	20.00	ng	0.00
21) Naphthalene-d8	8.96	136	332812	20.00	ng	0.00
38) Acenaphthene-d10	11.14	164	159755	20.00	ng	0.00
63) Phenanthrene-d10	12.98	188	311350	20.00	ng	0.00
75) Chrysene-d12	16.29	240	337558	20.00	ng	0.00
86) Perylene-d12	17.99	264	373259	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	415383	95.55	ng	0.01
7) Phenol-d6	6.92	99	549318	94.61	ng	0.00
23) Nitrobenzene-d5	8.07	82	298905	64.66	ng	0.00
41) 2,4,6-Tribromophenol	12.11	330	152389	97.32	ng	-0.01
44) 2-Fluorobiphenyl	10.31	172	665856	57.08	ng	0.00
78) Terphenyl-d14	14.98	244	753860	53.19	ng	0.00
Target Compounds						Qvalue
10) Phenol	6.94	94	19922	2.90	ng	# 71
31) Naphthalene	8.98	128	35290	2.23	ng	98
37) 2-Methylnaphthalene	9.85	142	42281	3.98	ng	97
70) Phenanthrene	13.02	178	116061	6.88	ng	98
74) Fluoranthene	14.49	202	73930	4.40	ng	98
77) Pyrene	14.78	202	72957	3.70	ng	97
80) Benzo(a)anthracene	16.26	228	56781m	3.14	ng	
82) Chrysene	16.30	228	179048m	10.14	ng	
85) Indeno(1,2,3-cd)pyrene	19.51	276	44483	2.22	ng	# 100
87) Benzo(b)fluoranthene	17.54	252	108757m	5.38	ng	
89) Benzo(a)pyrene	17.92	252	43148	2.39	ng	# 80
91) Benzo(a,h,i)perylene	19.95	276	59807	3.18	ng	# 91

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

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