

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042915\
 Data File : BF078838.D
 Acq On : 30 Apr 2015 5:10
 Operator : TP/IZ
 Sample : G2021-07
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PS-18-2-5

Quant Time: Apr 30 06:28:06 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	71825	20.00	ng	0.00
21) Naphthalene-d8	8.96	136	312117	20.00	ng	0.00
38) Acenaphthene-d10	11.14	164	150216	20.00	ng	0.00
63) Phenanthrene-d10	12.98	188	291168	20.00	ng	0.00
75) Chrysene-d12	16.27	240	318023	20.00	ng	-0.01
86) Perylene-d12	18.00	264	348424	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.68	112	209737	53.63	ng	0.01
7) Phenol-d6	6.92	99	329126	63.02	ng	0.00
23) Nitrobenzene-d5	8.07	82	170482	39.33	ng	0.00
41) 2,4,6-Tribromophenol	12.13	330	139936	95.17	ng	0.00
44) 2-Fluorobiphenyl	10.31	172	526361	47.99	ng	0.00
78) Terphenyl-d14	14.98	244	747433	55.97	ng	0.00
Target Compounds						Qvalue
70) Phenanthrene	13.02	178	213336	13.53	ng	99
74) Fluoranthene	14.49	202	254327	16.18	ng	98
77) Pyrene	14.78	202	289559	15.58	ng	98
80) Benzo(a)anthracene	16.26	228	132561	7.77	ng	99
82) Chrysene	16.31	228	115714	6.96	ng	99
85) Indeno(1,2,3-cd)pyrene	19.50	276	71878	3.81	ng	# 100
87) Benzo(b)fluoranthene	17.54	252	142060m	7.53	ng	
89) Benzo(a)pyrene	17.93	252	97344	5.77	ng	# 97
91) Benzo(g,h,i)perylene	19.95	276	74069	4.22	ng	98

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

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