

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051415\
 Data File : BF079237.D
 Acq On : 14 May 2015 17:13
 Operator : TP/IZ
 Sample : G2215-11
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-39S

Quant Time: May 15 12:48:19 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri May 15 02:08:55 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.22	152	64558	20.00	ng	0.00
21) Naphthalene-d8	8.80	136	270101	20.00	ng	0.00
38) Acenaphthene-d10	10.97	164	132202	20.00	ng	0.00
63) Phenanthrene-d10	12.82	188	257655	20.00	ng	0.00
75) Chrysene-d12	16.17	240	268237	20.00	ng	-0.33
86) Perylene-d12	18.29	264	271646	20.00	ng	-0.59
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	320758m	85.53	ng	0.00
7) Phenol-d6	6.79	99	443497	89.46	ng	0.00
23) Nitrobenzene-d5	7.91	82	233287	49.64	ng	-0.01
41) 2,4,6-Tribromophenol	11.95	330	123459	86.32	ng	0.00
44) 2-Fluorobiphenyl	10.15	172	445971	47.00	ng	0.00
78) Terphenyl-d14	14.82	244	475864m	42.47	ng	-0.10
Target Compounds						Qvalue
49) Dimethylphthalate	10.65	163	74830	7.71	ng	97
70) Phenanthrene	12.85	178	43608	3.03	ng	98
74) Fluoranthene	14.33	202	67694	4.51	ng	93
77) Pyrene	14.61	202	64793m	4.19	ng	
80) Benzo(a)anthracene	16.16	228	34810	2.37	ng	98
82) Chrysene	16.21	228	31813	2.25	ng	98
87) Benzo(b)fluoranthene	17.71	252	43055m	2.90	ng	
89) Benzo(a)pyrene	18.21	252	30504	2.23	ng	99

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

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