

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051215\
 Data File : BF079160.D
 Acq On : 12 May 2015 16:24
 Operator : TP/IZ
 Sample : G2155-10
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SB-43S

Quant Time: May 13 01:34:56 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 13 00:53:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.28	152	64455	20.00	ng	0.00
21) Naphthalene-d8	8.86	136	283279	20.00	ng	0.00
38) Acenaphthene-d10	11.03	164	135678	20.00	ng	-0.01
63) Phenanthrene-d10	12.88	188	263458	20.00	ng	0.00
75) Chrysene-d12	16.19	240	275564	20.00	ng	0.00
86) Perylene-d12	18.01	264	298496	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	296481	79.18	ng	0.01
7) Phenol-d6	6.84	99	400300	80.88	ng	0.00
23) Nitrobenzene-d5	7.98	82	223582	45.36	ng	0.00
41) 2,4,6-Tribromophenol	12.01	330	113031	77.01	ng	0.00
44) 2-Fluorobiphenyl	10.20	172	442921	45.48	ng	0.00
78) Terphenyl-d14	14.87	244	490605	42.62	ng	-0.01
Target Compounds						
49) Dimethylphthalate	10.71	163	22688	2.28	ng	Qvalue # 97
70) Phenanthrene	12.90	178	130436	8.85	ng	99
74) Fluoranthene	14.39	202	271630	17.69	ng	92
77) Pyrene	14.66	202	264808	16.68	ng	98
80) Benzo(a)anthracene	16.17	228	135135	8.96	ng	95
82) Chrysene	16.22	228	135514	9.34	ng	97
85) Indeno(1,2,3-cd)pyrene	19.47	276	103621	5.60	ng	# 100
87) Benzo(b)fluoranthene	17.54	252	178935m	10.95	ng	
88) Benzo(k)fluoranthene	17.55	252	74867m	4.73	ng	
89) Benzo(a)pyrene	17.94	252	137499	9.14	ng	# 97
91) Benzo(g,h,i)perylene	19.91	276	106021	6.62	ng	96

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

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