

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078592.D
 Acq On : 17 Apr 2015 5:38
 Operator : TP/IZ
 Sample : G1828-11DL 5X
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB02CDL

Quant Time: Apr 17 05:43:47 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 17 03:45:36 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	25037	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	109452	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	56814	20.00	ng	0.00
63) Phenanthrene-d10	12.24	188	117492	20.00	ng	0.00
75) Chrysene-d12	15.49	240	129440	20.00	ng	-0.01
86) Perylene-d12	17.15	264	133842	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	4.94	112	1323	0.87	ng	0.02
7) Phenol-d6	6.31	99	5303	2.79	ng	0.01
23) Nitrobenzene-d5	7.39	82	3895	2.16	ng	0.00
41) 2,4,6-Tribromophenol	11.39	330	1143	2.43	ng	0.00
44) 2-Fluorobiphenyl	9.62	172	7304	1.84	ng	0.00
78) Terphenyl-d14	14.23	244	36724	6.41	ng	0.00
Target Compounds						Qvalue
57) Fluorene	11.10	166	9553	2.32	ng	94
70) Phenanthrene	12.27	178	129935	19.12	ng	99
71) Anthracene	12.33	178	34536	5.16	ng	98
72) Carbazole	12.55	167	22503	3.59	ng	99
74) Fluoranthene	13.73	202	151340	21.12	ng	97
77) Pyrene	14.00	202	131155	16.14	ng	# 93
80) Benzo(a)anthracene	15.47	228	82297	10.69	ng	98
82) Chrysene	15.52	228	65165m	9.01	ng	
85) Indeno(1,2,3-cd)pyrene	18.32	276	26494	3.23	ng	# 66
87) Benzo(b)fluoranthene	16.73	252	81682m	9.87	ng	
88) Benzo(k)fluoranthene	16.75	252	17890m	2.43	ng	
89) Benzo(a)pyrene	17.10	252	55890	7.74	ng	98
91) Benzo(g,h,i)perylene	18.65	276	21837	3.01	ng	97

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

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