

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078593.D
 Acq On : 17 Apr 2015 6:07
 Operator : TP/IZ
 Sample : G1828-13DL 5X
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB01BDL

Quant Time: Apr 17 05:54:00 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	24091	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	98581	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	52803	20.00	ng	0.00
63) Phenanthrene-d10	12.24	188	110562	20.00	ng	0.00
75) Chrysene-d12	15.49	240	126037	20.00	ng	-0.01
86) Perylene-d12	17.15	264	129234	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	4.91	112	30518	20.89	ng	0.00
7) Phenol-d6	6.30	99	41375	22.60	ng	0.00
23) Nitrobenzene-d5	7.39	82	21722	13.36	ng	0.00
41) 2,4,6-Tribromophenol	11.39	330	9464	21.61	ng	0.00
44) 2-Fluorobiphenyl	9.62	172	34596	9.37	ng	0.00
78) Terphenyl-d14	14.23	244	50496	9.05	ng	0.00
Target Compounds						Qvalue
57) Fluorene	11.10	166	8512	2.23	ng	93
70) Phenanthrene	12.27	178	97157	15.19	ng	98
71) Anthracene	12.33	178	33725	5.35	ng	98
74) Fluoranthene	13.73	202	134704	19.97	ng	98
77) Pyrene	14.00	202	122000	15.42	ng	# 95
80) Benzo(a)anthracene	15.47	228	63882	8.52	ng	97
82) Chrysene	15.52	228	50892	7.22	ng	97
85) Indeno(1,2,3-cd)pyrene	18.32	276	27190	3.40	ng	# 82
87) Benzo(b)fluoranthene	16.73	252	65953m	8.26	ng	
88) Benzo(k)fluoranthene	16.77	252	15171m	2.14	ng	
89) Benzo(a)pyrene	17.09	252	46315	6.64	ng	# 90
91) Benzo(g,h,i)perylene	18.65	276	21603	3.09	ng	99

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

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