

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078571.D
 Acq On : 16 Apr 2015 17:54
 Operator : TP/IZ
 Sample : G1828-08RE
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB02BRE

Quant Time: Apr 17 02:00:04 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 00:52:22 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.68	152	22292	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	95460	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	48162	20.00	ng	0.00
63) Phenanthrene-d10	12.25	188	94888	20.00	ng	0.01
75) Chrysene-d12	15.51	240	107540	20.00	ng	0.02
86) Perylene-d12	17.27	264	116778	20.00	ng	0.08
System Monitoring Compounds						
5) 2-Fluorophenol	4.94	112	31201m	23.08	ng	0.03
7) Phenol-d6	6.31	99	63366	37.40	ng	0.01
23) Nitrobenzene-d5	7.39	82	43303	27.50	ng	0.00
41) 2,4,6-Tribromophenol	11.41	330	22696	56.82	ng	0.01
44) 2-Fluorobiphenyl	9.62	172	95339	28.30	ng	0.00
78) Terphenyl-d14	14.24	244	346901	72.89	ng	0.01
Target Compounds						Qvalue
48) Acenaphthylene	10.25	152	12038	2.41	ng	# 87
70) Phenanthrene	12.27	178	36069	6.57	ng	# 95
71) Anthracene	12.34	178	13850	2.56	ng	# 76
74) Fluoranthene	13.74	202	105760	18.27	ng	# 98
77) Pyrene	14.01	202	130240	19.29	ng	# 94
80) Benzo(a)anthracene	15.50	228	140907	22.02	ng	# 96
82) Chrysene	15.54	228	98016m	16.30	ng	
85) Indeno(1,2,3-cd)pyrene	18.51	276	216156m	31.69	ng	
87) Benzo(b)fluoranthene	16.82	252	247430m	34.28	ng	
88) Benzo(k)fluoranthene	16.85	252	82285m	12.83	ng	
89) Benzo(a)pyrene	17.20	252	221896	35.23	ng	# 92
90) Dibenzo(a,h)anthracene	18.53	278	39774m	6.31	ng	
91) Benzo(g,h,i)perylene	18.51	276	212343m	33.58	ng	

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

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