

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF041615\
 Data File : BF078591.D
 Acq On : 17 Apr 2015 5:09
 Operator : TP/IZ
 Sample : G1828-04DL 5X
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HS-SB06BDL

Quant Time: Apr 17 05:37:06 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	24465	20.00	ng	0.00
21) Naphthalene-d8	8.27	136	100094	20.00	ng	0.00
38) Acenaphthene-d10	10.42	164	53912	20.00	ng	0.00
63) Phenanthrene-d10	12.24	188	113242	20.00	ng	0.00
75) Chrysene-d12	15.49	240	128123	20.00	ng	-0.01
86) Perylene-d12	17.12	264	130704	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	4.91	112	25831	17.41	ng	0.00
7) Phenol-d6	6.30	99	37326	20.07	ng	0.00
23) Nitrobenzene-d5	7.39	82	20169	12.21	ng	0.00
41) 2,4,6-Tribromophenol	11.39	330	8683	19.42	ng	0.00
44) 2-Fluorobiphenyl	9.62	172	48293	12.80	ng	0.00
78) Terphenyl-d14	14.23	244	62705	11.06	ng	0.00
Target Compounds						Qvalue
51) Acenaphthene	10.46	154	6955	2.20	ng	98
57) Fluorene	11.10	166	8406	2.15	ng	99
70) Phenanthrene	12.27	178	111618	17.04	ng	99
71) Anthracene	12.33	178	21334	3.31	ng	97
74) Fluoranthene	13.73	202	187722	27.17	ng	96
77) Pyrene	14.00	202	154843	19.25	ng	# 95
80) Benzo(a)anthracene	15.47	228	70410	9.24	ng	97
82) Chrysene	15.52	228	65839	9.19	ng	97
85) Indeno(1,2,3-cd)pyrene	18.26	276	42379	5.21	ng	# 78
87) Benzo(b)fluoranthene	16.72	252	92580m	11.46	ng	
88) Benzo(k)fluoranthene	16.74	252	28756m	4.01	ng	
89) Benzo(a)pyrene	17.06	252	58038	8.23	ng	# 96
91) Benzo(g,h,i)perylene	18.58	276	36569	5.17	ng	95

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

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