

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111715\
 Data File : BF083004.D
 Acq On : 18 Nov 2015 9:49
 Operator : UM/IZ
 Sample : G4443-04
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 20151110MGP206BRV40

Quant Time: Nov 18 15:41:51 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 18 13:32:32 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.76	152	160467	20.00	ng	0.02
21) Naphthalene-d8	8.08	136	338328	20.00	ng	0.06
38) Acenaphthene-d10	9.79	164	406730	20.00	ng	0.01
63) Phenanthrene-d10	11.28	188	800100	20.00	ng	0.02
75) Chrysene-d12	13.91	240	726380	20.00	ng	0.01
86) Perylene-d12	15.30	264	543673	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.37	112	377490	36.60	ng	0.03
7) Phenol-d6	6.41	99	402411	29.35	ng	0.01
23) Nitrobenzene-d5	7.32	82	984043	126.56	ng	0.01
41) 2,4,6-Tribromophenol	10.58	330	269993	57.89	ng	0.01
44) 2-Fluorobiphenyl	9.13	172	2222159	78.30	ng	0.02
78) Terphenyl-d14	12.87	244	1852039	60.47	ng	0.02
Target Compounds						Qvalue
10) Phenol	6.42	94	126923	8.16	ng	80
20) 3+4-Methylphenols	7.18	107	37665	2.99	ng	# 61
27) 2,4-Dimethylphenol	7.74	122	48729	8.17	ng	89
31) Naphthalene	8.14	128	29766285	1594.62	ng	# 79
37) 2-Methylnaphthalene	8.78	142	7299351	604.48	ng	98
45) 1,1'-Biphenyl	9.22	154	1339719	38.39	ng	96
48) Acenaphthylene	9.66	152	4594197	107.68	ng	# 92
49) Dimethylphthalate	9.52	163	109362	3.28	ng	98
51) Acenaphthene	9.84	154	1810983	66.97	ng	99
54) Dibenzofuran	10.00	168	452829	12.42	ng	# 87
57) Fluorene	10.35	166	2263504	76.66	ng	99
70) Phenanthrene	11.30	178	2607447	56.12	ng	99
71) Anthracene	11.34	178	636932	13.01	ng	97
72) Carbazole	11.50	167	843073	19.68	ng	97
74) Fluoranthene	12.48	202	249317	5.00	ng	98
77) Pyrene	12.71	202	363303	7.28	ng	97

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

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