

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF111315\
 Data File : BF082904.D
 Acq On : 13 Nov 2015 18:30
 Operator : UM/IZ
 Sample : G4406-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PURGEWATER

Quant Time: Nov 16 09:36:18 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF110715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Nov 07 02:32:49 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	193899	20.00	ng	-0.06
21) Naphthalene-d8	8.09	136	766904	20.00	ng	-0.06
38) Acenaphthene-d10	9.84	164	349256	20.00	ng	-0.06
63) Phenanthrene-d10	11.32	188	692526	20.00	ng	-0.06
75) Chrysene-d12	13.95	240	531347	20.00	ng	-0.07
86) Perylene-d12	15.37	264	451956	20.00	ng	-0.09
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	670042	53.77	ng	-0.06
7) Phenol-d6	6.43	99	542107	32.72	ng	-0.05
23) Nitrobenzene-d5	7.37	82	1355072	76.88	ng	-0.04
41) 2,4,6-Tribromophenol	10.64	330	493158	123.15	ng	-0.04
44) 2-Fluorobiphenyl	9.17	172	2110890	86.62	ng	-0.05
78) Terphenyl-d14	12.91	244	1188085	53.03	ng	-0.06
Target Compounds						Qvalue
51) Acenaphthene	9.87	154	222829	9.60	ng	98
54) Dibenzofuran	10.04	168	189017	6.04	ng	90
57) Fluorene	10.38	166	127736	5.04	ng	97
59) Diethylphthalate	10.26	149	251692	9.01	ng	96
70) Phenanthrene	11.35	178	120441	2.99	ng	95
72) Carbazole	11.55	167	229951	6.20	ng	96

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

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