

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071315\
 Data File : BF080437.D
 Acq On : 13 Jul 2015 20:32
 Operator : TP/UM
 Sample : G2945-09
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PO-002

Quant Time: Jul 14 08:14:53 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 11 05:35:30 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.14	152	22822	20.00	ng	-0.01
21) Naphthalene-d8	8.42	136	90705	20.00	ng	-0.01
38) Acenaphthene-d10	10.18	164	49057	20.00	ng	-0.01
63) Phenanthrene-d10	11.66	188	95232	20.00	ng	-0.02
75) Chrysene-d12	14.48	240	106980	20.00	ng	-0.06
86) Perylene-d12	16.20	264	96122	20.00	ng	-0.08
System Monitoring Compounds						
5) 2-Fluorophenol	5.77	112	50415	37.86	ng	-0.02
7) Phenol-d6	6.78	99	39913	22.49	ng	-0.01
23) Nitrobenzene-d5	7.70	82	107541	67.62	ng	-0.01
41) 2,4,6-Tribromophenol	10.97	330	39288	82.06	ng	-0.02
44) 2-Fluorobiphenyl	9.49	172	246468	67.67	ng	-0.01
78) Terphenyl-d14	13.29	244	317437	64.44	ng	-0.03
Target Compounds						
10) Phenol	6.80	94	313717	160.00	ng	Qvalue 89
20) 3+4-Methylphenols	7.55	107	15563	8.88	ng	# 23
32) Benzoic acid	8.17	122	24570	26.29	ng	98
49) Dimethylphthalate	9.89	163	34087	9.19	ng	97

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

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