

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF090215\
 Data File : BF081346.D
 Acq On : 2 Sep 2015 13:55
 Operator : UM/IZ
 Sample : G3461-01
 Misc : Double Internal standard added
 ALS Vial : 10 Sample Multiplier: 2

Instrument :
 BNA_F
 ClientSampleId :
 LAR9837

Quant Time: Sep 02 23:47:11 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 22:52:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.39	152	110145	40.00	ng	0.00
21) Naphthalene-d8	7.68	136	463866	40.00	ng	0.00
38) Acenaphthene-d10	9.42	164	202192	40.00	ng	0.00
63) Phenanthrene-d10	10.88	188	390979	40.00	ng	0.00
75) Chrysene-d12	13.49	240	308005	40.00	ng	-0.01
86) Perylene-d12	14.80	264	212016	40.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	4.95	112	448235	126.28	ng	0.02
7) Phenol-d6	6.08	99	587939	125.13	ng	0.02
23) Nitrobenzene-d5	6.97	82	468574	97.47	ng	0.00
41) 2,4,6-Tribromophenol	10.20	330	129582	143.95	ng	0.00
44) 2-Fluorobiphenyl	8.76	172	784179	99.40	ng	0.00
78) Terphenyl-d14	12.47	244	590434	89.25	ng	0.00
Target Compounds						
10) Phenol	6.10	94	2381739	437.12	ng	# 63
17) 2-Methylphenol	6.68	107	8488	2.72	ng	# 81
20) 3+4-Methylphenols	6.84	107	27741	6.59	ng	85

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

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