

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG081616\
 Data File : BG023734.D
 Acq On : 16 Aug 2016 19:18
 Operator : UM/SJ
 Sample : PB92660BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB92660BL

Quant Time: Aug 17 01:27:41 2016
 Quant Method : Z:\HPCHEM1\BNA_G\Methods\8270-BG080116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 19:22:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.10	152	93127	20.00	ng	-0.02
21) Naphthalene-d8	10.94	136	391221	20.00	ng	-0.01
38) Acenaphthene-d10	14.78	164	264484	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	705079	20.00	ng	0.00
75) Chrysene-d12	21.87	240	842919	20.00	ng	0.00
86) Perylene-d12	25.28	264	816874	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	741386	134.01	ng	-0.02
7) Phenol-d6	7.26	99	1086686	126.11	ng	-0.01
23) Nitrobenzene-d5	9.28	82	723249	91.91	ng	-0.01
41) 2,4,6-Tribromophenol	16.28	330	535626	138.46	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	1430113	91.20	ng	0.00
78) Terphenyl-d14	20.16	244	2585271	99.17	ng	0.00

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

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