

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG092216\
 Data File : BG024076.D
 Acq On : 22 Sep 2016 20:22
 Operator : UM/SJ
 Sample : PB93511BL
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB93511BL

Quant Time: Sep 23 01:55:46 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG083116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 19:01:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	53156	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	231287	20.00	ng	0.00
38) Acenaphthene-d10	14.73	164	187492	20.00	ng	0.00
63) Phenanthrene-d10	17.49	188	389356	20.00	ng	0.00
75) Chrysene-d12	21.80	240	386910	20.00	ng	0.00
86) Perylene-d12	25.15	264	355238	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	266244	87.94	ng	0.00
7) Phenol-d6	7.22	99	402361	86.33	ng	0.00
23) Nitrobenzene-d5	9.23	82	470424	90.80	ng	0.00
41) 2,4,6-Tribromophenol	16.23	330	200061	59.22	ng	0.00
44) 2-Fluorobiphenyl	13.34	172	947605	72.46	ng	0.00
78) Terphenyl-d14	20.10	244	1405321	82.72	ng	0.00

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

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