

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG012615\
 Data File : BG015931.D
 Acq On : 27 Jan 2015 2:40
 Operator : TP/IZ
 Sample : G1139-09
 Misc : S
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121B2B

Quant Time: Jan 27 04:08:38 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG012015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 20 17:47:40 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.69	152	83504	20.00	ng	0.00
21) Naphthalene-d8	10.47	136	356072	20.00	ng	0.00
38) Acenaphthene-d10	14.33	164	292015	20.00	ng	0.00
63) Phenanthrene-d10	17.08	188	838594	20.00	ng	0.00
75) Chrysene-d12	21.26	240	1224780	20.00	ng	0.00
86) Perylene-d12	23.52	264	1165588	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.28	112	244051	43.52	ng	0.00
7) Phenol-d6	6.87	99	248710	27.52	ng	0.00
23) Nitrobenzene-d5	8.84	82	940704	76.15	ng	0.00
41) 2,4,6-Tribromophenol	15.83	330	490938	91.84	ng	0.00
44) 2-Fluorobiphenyl	12.95	172	1682896	90.42	ng	0.00
78) Terphenyl-d14	19.71	244	3010655	64.15	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	10.52	128	445622	24.45	ng	96
37) 2-Methylnaphthalene	12.14	142	549792	37.99	ng	97
45) 1,1'-Biphenyl	13.16	154	75685	4.11	ng	93
51) Acenaphthene	14.39	154	34290	2.31	ng	94
57) Fluorene	15.38	166	85104	3.85	ng	88
70) Phenanthrene	17.12	178	170912	4.06	ng	95
72) Carbazole	17.49	167	88002	2.42	ng	95

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

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