

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110416\
 Data File : BG024693.D
 Acq On : 5 Nov 2016 2:26
 Operator : UM/SJ
 Sample : H5531-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S22-0011

Quant Time: Nov 05 03:42:32 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG102516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 26 11:14:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.27	152	91517	20.00	ng	0.00
21) Naphthalene-d8	11.09	136	383854	20.00	ng	0.00
38) Acenaphthene-d10	14.88	164	316907	20.00	ng	0.00
63) Phenanthrene-d10	17.63	188	708076	20.00	ng	0.00
75) Chrysene-d12	21.92	240	971582	20.00	ng	0.00
86) Perylene-d12	25.35	264	1001864	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.80	112	412449	73.87	ng	0.00
7) Phenol-d6	7.40	99	386892	50.51	ng	0.00
23) Nitrobenzene-d5	9.45	82	688385	81.65	ng	0.00
41) 2,4,6-Tribromophenol	16.37	330	636725	134.49	ng	0.00
44) 2-Fluorobiphenyl	13.51	172	1446763	75.88	ng	0.00
78) Terphenyl-d14	20.21	244	2258669	69.47	ng	0.00
Target Compounds						
10) Phenol	7.43	94	22085	2.80	ng	Qvalue 88
22) Acetophenone	9.11	105	3818397	387.26	ng	# 95
27) 2,4-Dimethylphenol	10.24	122	155859	25.57	ng	99
32) Benzoic acid	10.36	122	12642	2.49	ng	# 70
49) Dimethylphthalate	14.32	163	106871	4.75	ng	# 88

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

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