

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110716\
 Data File : BG024727.D
 Acq On : 7 Nov 2016 13:20
 Operator : UM/SJ
 Sample : H5531-02DL 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S22-0013DL

Quant Time: Nov 07 15:28:06 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	8.26	152	93502	20.00	ng	0.00
21) Naphthalene-d8	11.09	136	372964	20.00	ng	-0.01
38) Acenaphthene-d10	14.88	164	298394	20.00	ng	-0.01
63) Phenanthrene-d10	17.62	188	677284	20.00	ng	0.00
75) Chrysene-d12	21.91	240	974260	20.00	ng	-0.02
86) Perylene-d12	25.34	264	1008223	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.79	112	110510	19.37	ng	0.00
7) Phenol-d6	7.39	99	102276	13.07	ng	0.00
23) Nitrobenzene-d5	9.44	82	165921	20.25	ng	-0.01
41) 2,4,6-Tribromophenol	16.36	330	156666	35.14	ng	0.00
44) 2-Fluorobiphenyl	13.50	172	401504	22.36	ng	-0.01
78) Terphenyl-d14	20.20	244	784106	24.05	ng	-0.01
Target Compounds						Qvalue
22) Acetophenone	9.08	105	304749	31.81	ng	# 88
31) Naphthalene	11.14	128	520535	27.98	ng	98
32) Benzoic acid	10.32	122	42741	8.67	ng	86
37) 2-Methylnaphthalene	12.73	142	144504	10.65	ng	95

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

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