

Data Path : Z:\HPCHEM1\BNA G\DATA\BG110416\
 Data File : BG024690.D
 Acq On : 5 Nov 2016 00:31
 Operator : UM/SJ
 Sample : H5531-02
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S22-0013

Manual Integrations
 APPROVED

sohil
 11/7/2016 4:36:48 PM

Quant Time: Nov 05 04:45:33 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	94204	20.00	ng	0.00
21) Naphthalene-d8	11.10	136	376076	20.00	ng	0.00
38) Acenaphthene-d10	14.89	164	307793	20.00	ng	0.00
63) Phenanthrene-d10	17.62	188	670674	20.00	ng	0.00
75) Chrysene-d12	21.92	240	971798	20.00	ng	-0.01
86) Perylene-d12	25.35	264	1003079	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.80	112	526740	91.64	ng	0.00
7) Phenol-d6	7.41	99	483445	61.32	ng	0.00
23) Nitrobenzene-d5	9.44	82	814554	98.61	ng	0.00
41) 2,4,6-Tribromophenol	16.37	330	700883	152.42	ng	0.00
44) 2-Fluorobiphenyl	13.51	172	1667169	90.03	ng	0.00
78) Terphenyl-d14	20.20	244	2910877	89.50	ng	-0.01
Target Compounds						
10) Phenol	7.44	94	19331	2.38	ng	# 75
22) Acetophenone	9.09	105	1387367	143.62	ng	# 92
27) 2,4-Dimethylphenol	10.24	122	111814m	18.73	ng	
31) Naphthalene	11.15	128	2302469	122.74	ng	97
32) Benzoic acid	10.41	122	246420	49.56	ng	90
37) 2-Methylnaphthalene	12.73	142	648769	47.40	ng	99
45) 1,1'-Biphenyl	13.72	154	62288	2.92	ng	91
49) Dimethylphthalate	14.32	163	109881	5.03	ng	# 94

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

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