

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122116\
 Data File : BG025369.D
 Acq On : 21 Dec 2016 22:41
 Operator : UM/SJ
 Sample : H6103-05
 Misc : LOD 8 PPM
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOD-WATER-05-QT1-2017

Manual Integrations
 APPROVED

sohil
 12/22/2016 7:04:35 PM

Quant Time: Dec 22 04:51:13 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 18:42:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.21	152	50388	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	209686	20.00	ng	0.00
38) Acenaphthene-d10	14.84	164	165072	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	395206	20.00	ng	0.00
75) Chrysene-d12	21.87	240	576343	20.00	ng	0.00
86) Perylene-d12	25.28	264	605170	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	419916	151.48	ng	0.00
7) Phenol-d6	7.37	99	581935	149.06	ng	0.00
23) Nitrobenzene-d5	9.40	82	441106	92.93	ng	0.00
41) 2,4,6-Tribromophenol	16.33	330	359859	135.41	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	958260	93.98	ng	0.00
78) Terphenyl-d14	20.16	244	1905304	87.65	ng	0.00
Target Compounds						
40) Hexachlorocyclopentadiene	13.01	237	3047	9.88	ng	Qvalue 94
53) 2,4-Dinitrophenol	15.02	184	2084m	6.21	ng	

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

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