

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121918\
 Data File : BG038751.D
 Acq On : 20 Dec 2018 6:57
 Operator : JU/SJ
 Sample : J6425-10
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-20181214

Quant Time: Dec 20 07:53:55 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 12 15:26:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.06	152	22348	20.00	ng	0.00
21) Naphthalene-d8	10.87	136	90559	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	62058	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	163252	20.00	ng	0.00
76) Chrysene-d12	21.72	240	174083	20.00	ng	0.00
87) Perylene-d12	25.01	264	178611	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	173844	137.97	ng	0.00
7) Phenol-d6	7.21	99	219008	126.61	ng	0.00
23) Nitrobenzene-d5	9.23	82	165123	93.15	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	117198	137.03	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	406871	89.94	ng	0.00
79) Terphenyl-d14	20.04	244	857567	107.21	ng	0.00
Target Compounds						
10) Phenol	7.24	94	7897	4.297	ng	85
20) 3+4-Methylphenols	8.82	107	7196	4.232	ng	93

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

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