

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA G\DATA\BG022218\
 Data File : BG032787.D
 Acq On : 22 Feb 2018 18:58
 Operator : SJ/JU
 Sample : J1637-03
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FRAC-TANK-A124

Quant Time: Feb 23 01:13:28 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 21 15:45:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.93	152	78274	20.00	ng	0.00
21) Naphthalene-d8	11.81	136	348832	20.00	ng	0.00
38) Acenaphthene-d10	15.54	164	155747	20.00	ng	-0.02
63) Phenanthrene-d10	18.27	188	405207	20.00	ng	0.00
75) Chrysene-d12	22.62	240	380640	20.00	ng	-0.02
86) Perylene-d12	26.43	264	77645	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.34	112	579413	118.43	ng	0.00
7) Phenol-d6	8.02	99	651158	95.48	ng	0.00
23) Nitrobenzene-d5	10.13	82	515938	100.58	ng	0.00
41) 2,4,6-Tribromophenol	17.00	330	290440	177.35	ng	0.00
44) 2-Fluorobiphenyl	14.17	172	1177949	109.08	ng	0.00
78) Terphenyl-d14	20.79	244	1569843	92.66	ng	0.00
Target Compounds						
10) Phenol	8.05	94	30914	4.497	ng	94
32) Benzoic acid	11.13	122	558861	110.290	ng	# 84
49) Dimethylphthalate	14.96	163	34029	2.573	ng	96

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

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