

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_G\DATA\BG010318\
 Data File : BG031905.D
 Acq On : 4 Jan 2018 00:45
 Operator : SJ/JU
 Sample : J1005-02
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 OWS-2

Manual Integrations
 APPROVED

Sohil
 1/4/2018 1:52:52 PM

Quant Time: Jan 04 08:18:52 2018
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 29 15:51:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.79	152	54490	20.00	ng	0.00
21) Naphthalene-d8	11.67	136	225089	20.00	ng	0.00
38) Acenaphthene-d10	15.42	164	151985	20.00	ng	-0.01
63) Phenanthrene-d10	18.15	188	377777	20.00	ng	-0.01
75) Chrysene-d12	22.49	240	427889	20.00	ng	-0.02
86) Perylene-d12	26.22	264	457805	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	6.21	112	423237	141.47	ng	0.00
7) Phenol-d6	7.89	99	512183	131.86	ng	0.00
23) Nitrobenzene-d5	9.99	82	324499	108.86	ng	0.00
41) 2,4,6-Tribromophenol	16.89	330	378286	163.12	ng	0.00
44) 2-Fluorobiphenyl	14.05	172	1145695	94.61	ng	0.00
78) Terphenyl-d14	20.68	244	1533191	81.86	ng	-0.01
Target Compounds						
9) Benzaldehyde	7.89	77	97533	41.702	ng	Qvalue 80
15) Benzyl Alcohol	9.05	79	1159853	397.777	ng	89
31) Naphthalene	11.73	128	37382	3.291	ng	100
32) Benzoic acid	10.90	122	73724	34.141	ng	# 76
58) 2,3,4,6-Tetrachlorophenol	16.02	232	13966m	3.500	ng	

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

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