

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG092118\
 Data File : BG036889.D
 Acq On : 22 Sep 2018 11:37
 Operator : JU/SJ
 Sample : J4993-05
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 NS-SLUDGE

Manual Integrations
 APPROVED

Sohil
 9/24/2018 9:29:58 AM

Quant Time: Sep 24 04:12:51 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 20 20:07:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	36207	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	158327	20.00	ng	0.00
38) Acenaphthene-d10	14.67	164	109037	20.00	ng	0.00
63) Phenanthrene-d10	17.42	188	283276	20.00	ng	0.00
75) Chrysene-d12	21.69	240	302550	20.00	ng	0.00
86) Perylene-d12	24.89	264	318769	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	193987	102.75	ng	0.00
7) Phenol-d6	7.20	99	267039	91.42	ng	-0.01
23) Nitrobenzene-d5	9.21	82	252436	85.38	ng	0.00
41) 2,4,6-Tribromophenol	16.16	330	149356	114.49	ng	-0.01
44) 2-Fluorobiphenyl	13.29	172	592269	80.90	ng	0.00
78) Terphenyl-d14	20.03	244	1058793	92.02	ng	0.00
Target Compounds						
10) Phenol	7.23	94	12001	3.981	ng	Qvalue # 75
20) 3+4-Methylphenols	8.82	107	10562m	3.616	ng	

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

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