

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121318\
 Data File : BG038565.D
 Acq On : 14 Dec 2018 12:28
 Operator : JU/SJ
 Sample : J6329-01
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 071-PIPE-1

Manual Integrations
 APPROVED

Sohil
 12/14/2018 5:09:24 PM

Quant Time: Dec 14 13:25:57 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	26326	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	106459	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	77021	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	177975	20.00	ng	0.00
76) Chrysene-d12	21.72	240	179876	20.00	ng	0.00
87) Perylene-d12	25.02	264	199030	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	171224	115.36	ng	0.00
7) Phenol-d6	7.21	99	228482	112.13	ng	0.00
23) Nitrobenzene-d5	9.23	82	143706	68.96	ng	0.00
42) 2,4,6-Tribromophenol	16.18	330	117002	110.22	ng	0.00
45) 2-Fluorobiphenyl	13.32	172	342742	61.05	ng	0.00
79) Terphenyl-d14	20.04	244	561562	67.94	ng	0.00
Target Compounds						
10) Phenol	7.24	94	16267	7.514	ng	95
50) Dimethylphthalate	14.14	163	32655	5.192	ng	98
75) Fluoranthene	19.49	202	43163	3.780	ng	98
78) Pyrene	19.86	202	52921	4.453	ng	95
81) Benzo(a)anthracene	21.70	228	25131m	2.293	ng	
83) Chrysene	21.77	228	22003	2.084	ng	97
88) Benzo(b)fluoranthene	23.96	252	26539	2.429	ng	# 80
90) Benzo(a)pyrene	24.86	252	24254	2.301	ng	# 92

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

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