

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG121318\
 Data File : BG038567.D
 Acq On : 14 Dec 2018 13:47
 Operator : JU/SJ
 Sample : J6319-01
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 055_BT-1-1

Manual Integrations
 APPROVED

Sohil
 12/14/2018 5:09:25 PM

Quant Time: Dec 14 15:33:57 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	29315	20.00	ng	0.00
21) Naphthalene-d8	10.88	136	120990	20.00	ng	0.00
39) Acenaphthene-d10	14.69	164	84423	20.00	ng	0.00
64) Phenanthrene-d10	17.44	188	191035	20.00	ng	0.00
76) Chrysene-d12	21.72	240	186518	20.00	ng	0.00
87) Perylene-d12	25.02	264	205752	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	120708	73.03	ng	0.00
7) Phenol-d6	7.22	99	168697	74.35	ng	0.00
23) Nitrobenzene-d5	9.24	82	111647	47.14	ng	0.00
42) 2,4,6-Tribromophenol	16.19	330	78722	67.66	ng	0.00
45) 2-Fluorobiphenyl	13.31	172	250677	40.73	ng	0.00
79) Terphenyl-d14	20.04	244	379543	44.29	ng	0.00
Target Compounds						
10) Phenol	7.24	94	10802	4.481	ng	72
38) 1-Methylnaphthalene	12.75	142	142174	34.451	ng	99
50) Dimethylphthalate	14.14	163	24221	3.513	ng	98
71) Phenanthrene	17.49	178	59111	5.967	ng	98
75) Fluoranthene	19.49	202	91345	7.452	ng	96
78) Pyrene	19.85	202	93906	7.621	ng	97
81) Benzo(a)anthracene	21.70	228	51457	4.527	ng	97
83) Chrysene	21.77	228	45528	4.159	ng	98
86) Indeno(1,2,3-cd)pyrene	28.79	276	44170	3.432	ng	# 96
88) Benzo(b)fluoranthene	23.96	252	74880	6.629	ng	99
89) Benzo(k)fluoranthene	24.02	252	26013m	2.312	ng	
90) Benzo(a)pyrene	24.85	252	60830m	5.582	ng	
92) Benzo(a,h,i)perylene	29.97	276	44000	4.115	ng	94

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

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