

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
 Data File : BG039216.D
 Acq On : 18 Jan 2019 15:16
 Operator : JU/SJ
 Sample : K1157-09
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-05-011519

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:11:32 PM

Quant Time: Jan 18 16:24:28 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG010919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 18 04:20:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	21751	20.00	ng	0.00
21) Naphthalene-d8	10.82	136	86639	20.00	ng	-0.01
39) Acenaphthene-d10	14.65	164	51547	20.00	ng	0.00
64) Phenanthrene-d10	17.40	188	128394	20.00	ng	0.00
76) Chrysene-d12	21.67	240	129919	20.00	ng	-0.01
87) Perylene-d12	24.93	264	145888	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	101536	88.55	ng	0.00
7) Phenol-d6	7.17	99	142288	86.23	ng	0.00
23) Nitrobenzene-d5	9.19	82	93736	59.20	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	70344	81.41	ng	-0.01
45) 2-Fluorobiphenyl	13.27	172	209501	55.62	ng	0.00
79) Terphenyl-d14	20.00	244	318164	45.30	ng	0.00
Target Compounds						
10) Phenol	7.20	94	8754	5.291	ng	90
50) Dimethylphthalate	14.10	163	18064	4.147	ng	# 95
71) Phenanthrene	17.44	178	98452	14.507	ng	97
72) Anthracene	17.53	178	19721	2.939	ng	99
75) Fluoranthene	19.45	202	169950	20.201	ng	97
78) Pyrene	19.81	202	145038	17.518	ng	99
81) Benzo(a)anthracene	21.65	228	101767	12.868	ng	98
83) Chrysene	21.72	228	93288m	12.402	ng	
86) Indeno(1,2,3-cd)pyrene	28.66	276	41393	4.605	ng	96
88) Benzo(b)fluoranthene	23.89	252	128072	15.921	ng	# 97
89) Benzo(k)fluoranthene	23.95	252	37407m	4.703	ng	
90) Benzo(a)pyrene	24.77	252	77957	10.026	ng	98
92) Benzo(a,h,i)perylene	29.86	276	38054	4.970	ng	# 96

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

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