

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG011819\
 Data File : BG039217.D
 Acq On : 18 Jan 2019 15:55
 Operator : JU/SJ
 Sample : K1157-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WC-06-011519

Manual Integrations
 APPROVED

Sohil
 1/21/2019 12:11:35 PM

Quant Time: Jan 18 22:04:21 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	19069	20.00	ng	0.00
21) Naphthalene-d8	10.83	136	81698	20.00	ng	0.00
39) Acenaphthene-d10	14.65	164	53047	20.00	ng	0.00
64) Phenanthrene-d10	17.39	188	109992	20.00	ng	0.00
76) Chrysene-d12	21.67	240	113959	20.00	ng	-0.02
87) Perylene-d12	24.93	264	135479	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	178002	177.08	ng	0.00
7) Phenol-d6	7.18	99	238567	164.91	ng	0.00
23) Nitrobenzene-d5	9.19	82	143169	95.89	ng	0.00
42) 2,4,6-Tribromophenol	16.14	330	121748	136.92	ng	0.00
45) 2-Fluorobiphenyl	13.27	172	439018	113.26	ng	0.00
79) Terphenyl-d14	20.00	244	559048	90.75	ng	0.00
Target Compounds						
10) Phenol	7.19	94	3789	2.612	ng	Qvalue # 41
50) Dimethylphthalate	14.10	163	33155	7.396	ng	98
71) Phenanthrene	17.44	178	17986	3.094	ng	96
75) Fluoranthene	19.45	202	39416	5.469	ng	98
78) Pyrene	19.81	202	38419	5.290	ng	98
81) Benzo(a)anthracene	21.65	228	19807	2.855	ng	98
83) Chrysene	21.72	228	21011	3.184	ng	95
88) Benzo(b)fluoranthene	23.90	252	27214m	3.643	ng	
90) Benzo(a)pyrene	24.77	252	20414	2.827	ng	# 98

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

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