

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
 Data File : BG043964.D
 Acq On : 7 Jan 2020 21:30
 Operator : JU
 Sample : L1037-01 5X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SPARTA-FINES

Quant Time: Jan 08 03:04:01 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG123019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 31 13:32:23 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.96	152	159226	20.00	ng	0.00
21) Naphthalene-d8	10.77	136	674760	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	413702	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	794201	20.00	ng	0.00
76) Chrysene-d12	21.59	240	645834	20.00	ng	0.00
87) Perylene-d12	24.72	264	728735	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	85587	9.87	ng	0.00
7) Phenol-d6	7.13	99	218328	18.01	ng	0.00
23) Nitrobenzene-d5	9.12	82	128207	10.16	ng	0.00
42) 2,4,6-Tribromophenol	16.08	330	15358	3.55	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	301377	13.20	ng	0.00
79) Terphenyl-d14	19.95	244	387575	5.70	ng	0.00
Target Compounds						Qvalue
71) Phenanthrene	17.38	178	115468	3.031	ng	93
75) Fluoranthene	19.39	202	219716	5.043	ng	95
78) Pyrene	19.74	202	197711	4.690	ng	92
81) Benzo(a)anthracene	21.57	228	90979	2.272	ng	95
83) Chrysene	21.63	228	99189	2.607	ng	94
88) Benzo(b)fluoranthene	23.73	252	133887	3.108	ng	# 87
90) Benzo(a)pyrene	24.57	252	90190	2.218	ng	# 92

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

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