

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG010720\
 Data File : BG043963.D
 Acq On : 7 Jan 2020 20:51
 Operator : JU
 Sample : L1017-01
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 TR-04-010620

Quant Time: Jan 08 03:01:01 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	7.97	152	166670	20.00	ng	0.00
21) Naphthalene-d8	10.76	136	705487	20.00	ng	0.00
39) Acenaphthene-d10	14.59	164	434233	20.00	ng	0.00
64) Phenanthrene-d10	17.34	188	834392	20.00	ng	0.00
76) Chrysene-d12	21.59	240	680344	20.00	ng	0.00
87) Perylene-d12	24.72	264	772838	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	505460	55.68	ng	0.00
7) Phenol-d6	7.13	99	783536	61.75	ng	0.00
23) Nitrobenzene-d5	9.12	82	494325	37.48	ng	0.00
42) 2,4,6-Tribromophenol	16.08	330	446170	98.19	ng	0.00
45) 2-Fluorobiphenyl	13.22	172	1436187	59.92	ng	0.00
79) Terphenyl-d14	19.95	244	2117406	71.91	ng	0.00
Target Compounds						
50) Dimethylphthalate	14.04	163	130937	4.247	ng	Qvalue # 97
71) Phenanthrene	17.38	178	110426	2.759	ng	92
75) Fluoranthene	19.39	202	225300	4.922	ng	91
78) Pyrene	19.75	202	246857	5.559	ng	93
81) Benzo(a)anthracene	21.57	228	119064	2.822	ng	96
83) Chrysene	21.63	228	120162	2.998	ng	93
88) Benzo(b)fluoranthene	23.73	252	123376	2.700	ng	# 96
90) Benzo(a)pyrene	24.57	252	99802	2.314	ng	98

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

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