

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG110419\
 Data File : BG043483.D
 Acq On : 4 Nov 2019 16:33
 Operator : JU
 Sample : K5146-15 2X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 CL-01-103019

Quant Time: Nov 04 17:11:40 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG102119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 04 15:22:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.01	152	52105	20.00	ng	0.00
21) Naphthalene-d8	10.81	136	200414	20.00	ng	0.00
39) Acenaphthene-d10	14.64	164	133147	20.00	ng	0.00
64) Phenanthrene-d10	17.38	188	295237	20.00	ng	0.00
76) Chrysene-d12	21.65	240	307153	20.00	ng	0.00
87) Perylene-d12	24.85	264	368299	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	120523	42.39	ng	0.00
7) Phenol-d6	7.20	99	169573	41.45	ng	0.00
23) Nitrobenzene-d5	9.18	82	96770	24.89	ng	0.00
42) 2,4,6-Tribromophenol	16.13	330	87456	34.52	ng	0.00
45) 2-Fluorobiphenyl	13.26	172	256353	26.60	ng	0.00
79) Terphenyl-d14	19.99	244	447700	27.35	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	14.09	163	86447	8.382	ng	97
71) Phenanthrene	17.43	178	49014	3.242	ng	99
75) Fluoranthene	19.44	202	98277	4.986	ng	97
78) Pyrene	19.80	202	86578	4.554	ng	98
81) Benzo(a)anthracene	21.63	228	48283	2.510	ng	97
83) Chrysene	21.70	228	46974	2.457	ng	95
88) Benzo(b)fluoranthene	23.83	252	63542	3.046	ng	# 88
90) Benzo(a)pyrene	24.69	252	47521	2.386	ng	# 92

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

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