

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG101519\
 Data File : BG043242.D
 Acq On : 16 Oct 2019 8:08
 Operator : JU
 Sample : K5326-01
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 IDW-SO-101019

Quant Time: Oct 16 09:11:27 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG092519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 09 01:11:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	55713	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	206417	20.00	ng	0.00
39) Acenaphthene-d10	14.68	164	145959	20.00	ng	0.00
64) Phenanthrene-d10	17.42	188	320157	20.00	ng	0.00
76) Chrysene-d12	21.70	240	342096	20.00	ng	0.00
87) Perylene-d12	24.92	264	408642	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.64	112	411713	131.88	ng	0.00
7) Phenol-d6	7.23	99	563286	125.30	ng	0.00
23) Nitrobenzene-d5	9.22	82	366252	86.24	ng	0.00
42) 2,4,6-Tribromophenol	16.17	330	345947	137.84	ng	0.00
45) 2-Fluorobiphenyl	13.30	172	972668	96.61	ng	0.00
79) Terphenyl-d14	20.03	244	1685094	104.96	ng	0.00
Target Compounds						
31) Naphthalene	10.91	128	164256	16.165	ng	97
37) 2-Methylnaphthalene	12.51	142	17447	2.251	ng	94
38) 1-Methylnaphthalene	12.73	142	18011	2.456	ng	94
71) Phenanthrene	17.46	178	42885	2.744	ng	95
73) Carbazole	17.83	167	120577	8.333	ng	95

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

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