

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG083019\
 Data File : BG042619.D
 Acq On : 31 Aug 2019 3:02
 Operator : HP/JU
 Sample : K4439-09
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S-4A-(5-7)

Quant Time: Aug 31 05:24:29 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG082219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 22 14:29:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.99	152	36293	20.00	ng	-0.01
21) Naphthalene-d8	10.82	136	139898	20.00	ng	0.00
39) Acenaphthene-d10	14.66	164	104661	20.00	ng	0.00
64) Phenanthrene-d10	17.41	188	233171	20.00	ng	-0.01
76) Chrysene-d12	21.68	240	246143	20.00	ng	-0.01
87) Perylene-d12	24.92	264	291155	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	169487	94.58	ng	0.00
7) Phenol-d6	7.16	99	227476	93.98	ng	-0.01
23) Nitrobenzene-d5	9.16	82	131574	64.99	ng	-0.01
42) 2,4,6-Tribromophenol	16.15	330	144644	97.23	ng	-0.01
45) 2-Fluorobiphenyl	13.28	172	410691	66.37	ng	0.00
79) Terphenyl-d14	20.02	244	704010	61.83	ng	0.00
Target Compounds						
10) Phenol	7.19	94	7207	2.928	ng	85
50) Dimethylphthalate	14.10	163	63590	8.421	ng	98
71) Phenanthrene	17.45	178	31847	2.750	ng	99
75) Fluoranthene	19.46	202	56260	3.980	ng	96
78) Pyrene	19.82	202	58915	3.904	ng	96
83) Chrysene	21.72	228	31397	2.174	ng	96
88) Benzo(b)fluoranthene	23.89	252	34851	2.177	ng	99

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

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