

Data Path : Z:\SVOASRV\HPCHEM1\BNA_G\DATA\BG040319\
 Data File : BG040334.D
 Acq On : 3 Apr 2019 20:26
 Operator : JU/SJ
 Sample : K2182-05
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8B(2-10)COMP

Quant Time: Apr 04 02:22:17 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG032719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 28 05:47:00 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	60678	20.00	ng	-0.02
21) Naphthalene-d8	10.86	136	240352	20.00	ng	-0.02
39) Acenaphthene-d10	14.68	164	150996	20.00	ng	-0.02
64) Phenanthrene-d10	17.42	188	368233	20.00	ng	-0.02
76) Chrysene-d12	21.70	240	408128	20.00	ng	-0.04
87) Perylene-d12	24.98	264	418534	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	372817	102.14	ng	-0.02
7) Phenol-d6	7.20	99	528895	104.85	ng	-0.02
23) Nitrobenzene-d5	9.22	82	309074	68.35	ng	-0.02
42) 2,4,6-Tribromophenol	16.17	330	189480	116.11	ng	-0.02
45) 2-Fluorobiphenyl	13.30	172	727347	71.38	ng	-0.02
79) Terphenyl-d14	20.03	244	1182257	65.17	ng	-0.02
Target Compounds						
10) Phenol	7.23	94	14997	2.739	ng	96
50) Dimethylphthalate	14.13	163	53799	4.552	ng	97
75) Fluoranthene	19.47	202	152426	6.651	ng	95
78) Pyrene	19.83	202	169042	6.298	ng	97
81) Benzo(a)anthracene	21.68	228	71041	2.826	ng	97
83) Chrysene	21.74	228	96569	3.997	ng	98
88) Benzo(b)fluoranthene	23.92	252	140598	5.487	ng	100

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

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