

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060619\
 Data File : BG041092.D
 Acq On : 6 Jun 2019 15:36
 Operator : HP/JU
 Sample : K3076-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 FESB-07(5-5.5)

Quant Time: Jun 06 23:46:56 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG052919.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed May 29 18:39:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.11	152	53069	20.00	ng	-0.02
21) Naphthalene-d8	10.93	136	203420	20.00	ng	-0.02
39) Acenaphthene-d10	14.74	164	135007	20.00	ng	-0.02
64) Phenanthrene-d10	17.49	188	336682	20.00	ng	-0.02
76) Chrysene-d12	21.76	240	352914	20.00	ng	-0.02
87) Perylene-d12	25.10	264	388824	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.65	112	353562	120.14	ng	-0.01
7) Phenol-d6	7.26	99	514401	113.17	ng	-0.01
23) Nitrobenzene-d5	9.29	82	351407	76.69	ng	-0.02
42) 2,4,6-Tribromophenol	16.23	330	246890	123.18	ng	-0.02
45) 2-Fluorobiphenyl	13.37	172	690882	73.70	ng	-0.02
79) Terphenyl-d14	20.08	244	915694	55.07	ng	-0.02
Target Compounds						
50) Dimethylphthalate	14.18	163	128174	11.215	ng	99
75) Fluoranthene	19.53	202	88602	4.090	ng	97
78) Pyrene	19.89	202	85187	3.713	ng	96
83) Chrysene	21.81	228	50786	2.259	ng	96
88) Benzo(b)fluoranthene	24.03	252	86528	3.669	ng	# 91
90) Benzo(a)pyrene	24.94	252	56597	2.515	ng	99
92) Benzo(g,h,i)perylene	30.12	276	49360	2.260	ng	# 94

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

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