

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG061919\
 Data File : BG041328.D
 Acq On : 19 Jun 2019 15:04
 Operator : HP/JU
 Sample : K3154-01 2X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 JC-02-061419-A

Quant Time: Jun 19 16:35:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG061719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 17 14:58:02 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	38700	20.00	ng	0.00
21) Naphthalene-d8	10.90	136	156655	20.00	ng	0.00
39) Acenaphthene-d10	14.71	164	106393	20.00	ng	0.00
64) Phenanthrene-d10	17.46	188	249968	20.00	ng	0.00
76) Chrysene-d12	21.74	240	256635	20.00	ng	0.00
87) Perylene-d12	25.05	264	283718	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.61	112	104310	49.52	ng	0.00
7) Phenol-d6	7.23	99	149748	49.05	ng	0.00
23) Nitrobenzene-d5	9.26	82	106890	28.70	ng	0.00
42) 2,4,6-Tribromophenol	16.20	330	69862	42.77	ng	0.00
45) 2-Fluorobiphenyl	13.33	172	230595	29.46	ng	0.00
79) Terphenyl-d14	20.06	244	372151	28.74	ng	0.00
Target Compounds						
71) Phenanthrene	17.51	178	31339	2.351	ng	98
75) Fluoranthene	19.51	202	89620	5.060	ng	99
78) Pyrene	19.87	202	86799	5.015	ng	97
81) Benzo(a)anthracene	21.72	228	51979	2.991	ng	96
83) Chrysene	21.78	228	51994	3.111	ng	96
88) Benzo(b)fluoranthene	23.99	252	59108	3.362	ng	# 89
90) Benzo(a)pyrene	24.89	252	46529	2.802	ng	92

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

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