

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF042319\
 Data File : BF113922.D
 Acq On : 23 Apr 2019 16:07
 Operator : JU/SJ
 Sample : K2201-09 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-01-041819-E

Quant Time: Apr 23 16:56:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF042219.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 23 03:32:35 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	373066	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	1309836	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	601713	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	1130638	20.00	ng	0.00
76) Chrysene-d12	14.06	240	1096618	20.00	ng	0.00
87) Perylene-d12	15.56	264	911554	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	482961	22.14	ng	0.00
7) Phenol-d6	6.52	99	620164	22.66	ng	0.00
23) Nitrobenzene-d5	7.45	82	368827	17.39	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	161456	22.03	ng	0.00
45) 2-Fluorobiphenyl	9.25	172	689723	17.93	ng	0.00
79) Terphenyl-d14	13.00	244	841823	15.74	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.64	163	121165	2.784	ng	99
71) Phenanthrene	11.45	178	181613	3.196	ng	97
75) Fluoranthene	12.63	202	359519	5.799	ng	99
78) Pyrene	12.86	202	359132	4.683	ng	97
81) Benzo(a)anthracene	14.05	228	208750	3.231	ng	95
83) Chrysene	14.09	228	197646	3.141	ng	97
88) Benzo(b)fluoranthene	15.12	252	198428	3.441	ng	# 90
90) Benzo(a)pyrene	15.49	252	131642	2.639	ng	# 92

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

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