

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062219\
 Data File : BF115127.D
 Acq On : 22 Jun 2019 18:10
 Operator : HP/JU
 Sample : K3421-02
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SAMPLE-1

Quant Time: Jun 24 05:37:26 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF062119.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 24 04:21:29 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.61	152	297947	20.00	ng	0.00
21) Naphthalene-d8	7.89	136	1126536	20.00	ng	-0.01
39) Acenaphthene-d10	9.64	164	551875	20.00	ng	-0.01
64) Phenanthrene-d10	11.11	188	1054296	20.00	ng	-0.01
76) Chrysene-d12	13.72	240	629846	20.00	ng	-0.02
87) Perylene-d12	15.08	264	660741	20.00	ng	-0.03
System Monitoring Compounds						
5) 2-Fluorophenol	5.21	112	1596228	82.67	ng	0.00
7) Phenol-d6	6.25	99	1986248	84.76	ng	0.00
23) Nitrobenzene-d5	7.17	82	1169850	63.88	ng	-0.01
42) 2,4,6-Tribromophenol	10.41	330	419967	72.16	ng	-0.02
45) 2-Fluorobiphenyl	8.97	172	2098670	64.60	ng	-0.01
79) Terphenyl-d14	12.69	244	2043924	69.00	ng	-0.01
Target Compounds						
50) Dimethylphthalate	9.36	163	315404	7.768	ng	99
71) Phenanthrene	11.13	178	351363	6.190	ng	96
75) Fluoranthene	12.31	202	503713	8.894	ng	97
78) Pyrene	12.53	202	389836	8.054	ng	99
81) Benzo(a)anthracene	13.71	228	116393	2.575	ng	99
83) Chrysene	13.75	228	107799	2.604	ng	96
88) Benzo(b)fluoranthene	14.71	252	96611m	2.343	ng	

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

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