

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102519\
 Data File : BF117539.D
 Acq On : 25 Oct 2019 16:39
 Operator : JU
 Sample : K5501-15
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 183-82-(0-2.1)

Quant Time: Oct 26 05:51:04 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 18 18:53:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	36159	20.00	ng	0.00
21) Naphthalene-d8	8.01	136	144299	20.00	ng	0.00
39) Acenaphthene-d10	9.76	164	70703	20.00	ng	0.00
64) Phenanthrene-d10	11.25	188	99137	20.00	ng	0.00
76) Chrysene-d12	13.89	240	70382	20.00	ng	0.00
87) Perylene-d12	15.33	264	74068	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.35	112	167567	68.94	ng	0.00
7) Phenol-d6	6.38	99	227985	69.84	ng	0.00
23) Nitrobenzene-d5	7.29	82	139763	47.82	ng	-0.01
42) 2,4,6-Tribromophenol	10.56	330	36938	60.38	ng	0.00
45) 2-Fluorobiphenyl	9.08	172	246484	49.12	ng	-0.01
79) Terphenyl-d14	12.83	244	172999	40.09	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.48	163	33396	6.626	ng	Qvalue 100
71) Phenanthrene	11.27	178	12025	2.089	ng	96
75) Fluoranthene	12.46	202	25113	4.358	ng	98
78) Pyrene	12.69	202	23041	3.811	ng	97
81) Benzo(a)anthracene	13.88	228	12853	2.707	ng	99
83) Chrysene	13.92	228	12935	2.780	ng	96
88) Benzo(b)fluoranthene	14.92	252	16868m	3.857	ng	
90) Benzo(a)pyrene	15.27	252	11034	2.775	ng	# 80

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

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