

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF090419\
 Data File : BF116826.D
 Acq On : 4 Sep 2019 20:03
 Operator : HP/JU
 Sample : K4668-01
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-090319

Quant Time: Sep 05 04:33:22 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF083019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 30 20:02:30 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.85	152	69134	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	271189	20.00	ng	0.00
39) Acenaphthene-d10	9.88	164	130902	20.00	ng	0.00
64) Phenanthrene-d10	11.36	188	186254	20.00	ng	0.00
76) Chrysene-d12	14.00	240	141174	20.00	ng	0.00
87) Perylene-d12	15.46	264	107263	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	401296	94.04	ng	0.00
7) Phenol-d6	6.48	99	544049	97.09	ng	0.00
23) Nitrobenzene-d5	7.40	82	330184	69.51	ng	0.00
42) 2,4,6-Tribromophenol	10.67	330	89805	102.62	ng	0.00
45) 2-Fluorobiphenyl	9.20	172	583652	75.21	ng	0.00
79) Terphenyl-d14	12.95	244	449820	58.94	ng	0.00
Target Compounds						
10) Phenol	6.49	94	13451	2.168	ng	94
50) Dimethylphthalate	9.59	163	69769	7.720	ng	98
71) Phenanthrene	11.39	178	27291	2.632	ng	99
75) Fluoranthene	12.57	202	39041	3.832	ng	99
78) Pyrene	12.80	202	34484	2.748	ng	97
88) Benzo(b)fluoranthene	15.03	252	14188m	2.188	ng	

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

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