

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122118\
 Data File : BF111691.D
 Acq On : 21 Dec 2018 17:00
 Operator : JU/SJ
 Sample : J6447-08
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 P001-SS025-1218-01

Quant Time: Dec 21 23:53:16 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.90	152	183156	20.00	ng	0.00
21) Naphthalene-d8	8.18	136	636422	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	271111	20.00	ng	0.00
64) Phenanthrene-d10	11.42	188	453580	20.00	ng	0.00
76) Chrysene-d12	14.06	240	342925	20.00	ng	0.01
87) Perylene-d12	15.56	264	255290	20.00	ng	0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	1069280	99.51	ng	0.01
7) Phenol-d6	6.53	99	1328484	97.15	ng	0.01
23) Nitrobenzene-d5	7.46	82	833381	76.22	ng	0.00
42) 2,4,6-Tribromophenol	10.72	330	313879	97.98	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	1434518	77.12	ng	0.00
79) Terphenyl-d14	13.01	244	1262351	69.81	ng	0.01
Target Compounds						
10) Phenol	6.55	94	90267	5.850	ng	99
40) 1,2,4,5-Tetrachlorobenzene	9.06	216	18790	2.109	ng	98
46) 1,1'-Biphenyl	9.36	154	54088	2.363	ng	91
50) Dimethylphthalate	9.65	163	144074	7.267	ng	100
68) Hexachlorobenzene	11.03	284	15656	2.496	ng	94

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

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