

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF122618\
 Data File : BF111740.D
 Acq On : 27 Dec 2018 3:23
 Operator : JU/SJ
 Sample : J6188-05 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-122118

Manual Integrations
 APPROVED

Sohil
 12/27/2018 8:22:43 AM

Quant Time: Dec 27 04:30:35 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	132316	20.00	ng	0.00
21) Naphthalene-d8	8.19	136	446452	20.00	ng	0.00
39) Acenaphthene-d10	9.94	164	204506	20.00	ng	0.00
64) Phenanthrene-d10	11.43	188	397791	20.00	ng	0.01
76) Chrysene-d12	14.07	240	278755	20.00	ng	0.02
87) Perylene-d12	15.55	264	222050	20.00	ng	0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	388391	50.03	ng	0.02
7) Phenol-d6	6.54	99	463108	46.88	ng	0.02
23) Nitrobenzene-d5	7.46	82	262052	34.17	ng	0.00
42) 2,4,6-Tribromophenol	10.73	330	119086	49.28	ng	0.00
45) 2-Fluorobiphenyl	9.26	172	447841	31.92	ng	0.00
79) Terphenyl-d14	13.01	244	466066	31.71	ng	0.01
Target Compounds						
10) Phenol	6.55	94	28551	2.561	ng	94
50) Dimethylphthalate	9.65	163	55741	3.727	ng	97
71) Phenanthrene	11.45	178	75066	3.558	ng	92
75) Fluoranthene	12.64	202	95509	4.471	ng	95
78) Pyrene	12.87	202	84856	4.311	ng	97
81) Benzo(a)anthracene	14.06	228	33603	2.112	ng	98
83) Chrysene	14.09	228	31423m	2.010	ng	
88) Benzo(b)fluoranthene	15.12	252	34729m	2.676	ng	

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

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