

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102218\
 Data File : BF110050.D
 Acq On : 22 Oct 2018 15:50
 Operator : JU/SJ
 Sample : J5197-03 2X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HD-02-101918

Manual Integrations
 APPROVED

Sohil
 10/23/2018 6:09:45 PM

Quant Time: Oct 23 13:15:06 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 16 14:58:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	199388	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	750301	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	294235	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	469115	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	354467	20.00	ng	-0.01
87) Perylene-d12	15.67	264	252725	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	494386	39.30	ng	-0.02
7) Phenol-d6	6.58	99	620302	39.71	ng	-0.02
23) Nitrobenzene-d5	7.53	82	358250	32.16	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	97022	38.09	ng	-0.02
45) 2-Fluorobiphenyl	9.33	172	639446	34.68	ng	-0.01
79) Terphenyl-d14	13.09	244	485924	28.79	ng	-0.01
Target Compounds						
10) Phenol	6.59	94	47914	2.841	ng	# 65
50) Dimethylphthalate	9.72	163	62370	2.994	ng	# 97
71) Phenanthrene	11.53	178	73434	3.024	ng	99
75) Fluoranthene	12.72	202	119684	4.933	ng	99
78) Pyrene	12.96	202	153772	5.906	ng	99
81) Benzo(a)anthracene	14.14	228	67710	3.243	ng	93
83) Chrysene	14.17	228	82435m	4.148	ng	
88) Benzo(b)fluoranthene	15.22	252	73613m	5.052	ng	
90) Benzo(a)pyrene	15.60	252	56780	4.237	ng	# 88
92) Benzo(g,h,i)perylene	17.71	276	34286	2.857	ng	# 90

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

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