

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101718\
 Data File : BF109930.D
 Acq On : 17 Oct 2018 19:54
 Operator : JU/SJ
 Sample : J5200-11 2X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-01-101618-B

Manual Integrations
 APPROVED

Sohil
 10/18/2018 2:04:29 PM

10/18/2018 2:04:29 PM

Quant Time: Oct 18 11:13:22 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	273267	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	1073628	20.00	ng	-0.01
39) Acenaphthene-d10	10.02	164	469851	20.00	ng	-0.01
64) Phenanthrene-d10	11.51	188	746024	20.00	ng	0.00
76) Chrysene-d12	14.15	240	613535	20.00	ng	-0.01
87) Perylene-d12	15.68	264	669183	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	644818	37.40	ng	-0.01
7) Phenol-d6	6.59	99	804138	37.56	ng	-0.02
23) Nitrobenzene-d5	7.53	82	452457	28.39	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	138477	34.04	ng	-0.01
45) 2-Fluorobiphenyl	9.33	172	720244	24.46	ng	0.00
79) Terphenyl-d14	13.09	244	537978	18.41	ng	-0.02
Target Compounds						Qvalue
50) Dimethylphthalate	9.72	163	92007	2.765	ng	98
71) Phenanthrene	11.53	178	116680	3.021	ng	99
75) Fluoranthene	12.72	202	133788	3.468	ng	97
78) Pyrene	12.95	202	127849	2.837	ng	99
81) Benzo(a)anthracene	14.14	228	73337	2.029	ng	98
88) Benzo(b)fluoranthene	15.22	252	93649m	2.427	ng	

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

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