

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF082118\
 Data File : BF108358.D
 Acq On : 22 Aug 2018 00:47
 Operator : JU/SJ
 Sample : J4271-13 10X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 HR-06-082018-C

Manual Integrations
 APPROVED

Sohil
 8/22/2018 4:03:18 PM

Quant Time: Aug 22 01:56:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF080818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 21 15:20:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	170106	20.00	ng	-0.01
21) Naphthalene-d8	8.29	136	635413	20.00	ng	-0.01
38) Acenaphthene-d10	10.05	164	287295	20.00	ng	-0.02
63) Phenanthrene-d10	11.54	188	470031	20.00	ng	-0.02
75) Chrysene-d12	14.19	240	383669	20.00	ng	-0.02
86) Perylene-d12	15.75	264	326216	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	86227	9.18	ng	-0.01
7) Phenol-d6	6.61	99	121105	9.70	ng	-0.02
23) Nitrobenzene-d5	7.56	82	74072	6.50	ng	-0.02
41) 2,4,6-Tribromophenol	10.84	330	21612	7.54	ng	-0.02
44) 2-Fluorobiphenyl	9.36	172	140890	7.87	ng	-0.02
78) Terphenyl-d14	13.12	244	106239	7.04	ng	-0.02
Target Compounds						
70) Phenanthrene	11.57	178	67796	3.058	ng	99
74) Fluoranthene	12.77	202	153307	6.336	ng	93
77) Pyrene	12.99	202	132704	5.463	ng	99
80) Benzo(a)anthracene	14.18	228	79508	3.611	ng	99
82) Chrysene	14.22	228	65163	3.228	ng	97
87) Benzo(b)fluoranthene	15.28	252	86115m	4.418	ng	
89) Benzo(a)pyrene	15.68	252	53188	3.047	ng	95

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

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