

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF080318\
 Data File : BF107950.D
 Acq On : 3 Aug 2018 20:54
 Operator : JU/SJ
 Sample : J4318-07 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BEEBEE-SITE-DEMO-3

Manual Integrations
 APPROVED

Sohil
 8/6/2018 5:05:46 PM

Quant Time: Aug 04 01:19:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 30 17:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	113978	20.00	ng	-0.01
21) Naphthalene-d8	8.24	136	551517	20.00	ng	-0.01
38) Acenaphthene-d10	9.99	164	227557	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	401982	20.00	ng	-0.01
75) Chrysene-d12	14.14	240	351860	20.00	ng	-0.01
86) Perylene-d12	15.68	264	344179	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.63	112	148824m	22.19	ng	0.04
7) Phenol-d6	6.62	99	241436	27.32	ng	0.01
23) Nitrobenzene-d5	7.54	82	114407	11.95	ng	0.00
41) 2,4,6-Tribromophenol	10.79	330	67453	21.82	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	394656	24.17	ng	-0.01
78) Terphenyl-d14	13.07	244	320052	19.72	ng	-0.01
Target Compounds						
70) Phenanthrene	11.51	178	73378	3.844	ng	98
74) Fluoranthene	12.71	202	152835	6.816	ng	97
77) Pyrene	12.94	202	118436	4.674	ng	97
80) Benzo(a)anthracene	14.13	228	64737	3.248	ng	# 95
82) Chrysene	14.17	228	73896	3.437	ng	# 96
87) Benzo(b)fluoranthene	15.22	252	71331m	4.202	ng	
88) Benzo(k)fluoranthene	15.24	252	53336m	2.559	ng	
89) Benzo(a)pyrene	15.61	252	44728	2.552	ng	# 87
91) Benzo(g,h,i)perylene	17.75	276	29947	2.025	ng	# 83

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

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