

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF072318\
 Data File : BF107588.D
 Acq On : 23 Jul 2018 15:18
 Operator : JU/SJ
 Sample : J4062-25
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TP-9

Manual Integrations
 APPROVED

Sohil
 7/24/2018 5:26:02 PM

Quant Time: Jul 24 14:30:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF072018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 20 16:29:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	319016	20.00	ng	0.00
21) Naphthalene-d8	8.25	136	1257719	20.00	ng	0.00
38) Acenaphthene-d10	10.02	164	529318	20.00	ng	0.00
63) Phenanthrene-d10	11.51	188	925616m	20.00	ng	0.00
75) Chrysene-d12	14.16	240	732568	20.00	ng	0.00
86) Perylene-d12	15.70	264	738096	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1518136	76.51	ng	0.00
7) Phenol-d6	6.61	99	1921330	80.64	ng	0.00
23) Nitrobenzene-d5	7.54	82	1213610	65.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.81	330	511374	84.96	ng	0.00
44) 2-Fluorobiphenyl	9.33	172	1974685	58.37	ng	0.00
78) Terphenyl-d14	13.10	244	1564400	54.67	ng	0.00
Target Compounds						
10) Phenol	6.62	94	73014	2.606	ng	# 72
49) Dimethylphthalate	9.72	163	104084	2.609	ng	# 99
70) Phenanthrene	11.54	178	148642	3.295	ng	99
74) Fluoranthene	12.73	202	245901	5.080	ng	96
77) Pyrene	12.96	202	226850	4.527	ng	97
80) Benzo(a)anthracene	14.15	228	107197	2.497	ng	96
82) Chrysene	14.19	228	109159	2.647	ng	97
87) Benzo(b)fluoranthene	15.24	252	125591m	3.109	ng	
89) Benzo(a)pyrene	15.63	252	82656	2.237	ng	95

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

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