

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\
 Data File : BF109233.D
 Acq On : 22 Sep 2018 20:35
 Operator : JU/SJ
 Sample : J4777-23 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-03-092018-C

Manual Integrations
 APPROVED

Sohil
 9/24/2018 11:39:01 AM

Quant Time: Sep 24 07:49:12 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	124373	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	475530	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	215529	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	364013	20.00	ng	0.00
75) Chrysene-d12	13.83	240	251521	20.00	ng	-0.01
86) Perylene-d12	15.23	264	238578	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	138500	18.34	ng	-0.01
7) Phenol-d6	6.33	99	179302	18.61	ng	-0.02
23) Nitrobenzene-d5	7.26	82	105069	13.04	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	36324	17.10	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	215498	15.08	ng	-0.01
78) Terphenyl-d14	12.79	244	156210	15.24	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.46	163	54043	3.447	ng	98
74) Fluoranthene	12.42	202	61260	3.154	ng	95
77) Pyrene	12.64	202	61117	3.390	ng	98
80) Benzo(a)anthracene	13.82	228	33768m	2.229	ng	
82) Chrysene	13.86	228	31130	2.073	ng	97
87) Benzo(b)fluoranthene	14.84	252	42095m	3.211	ng	
89) Benzo(a)pyrene	15.17	252	27699	2.345	ng	# 93

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

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