

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092218\
 Data File : BF109234.D
 Acq On : 22 Sep 2018 21:01
 Operator : JU/SJ
 Sample : J4777-25 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-04-092018-A

Manual Integrations
 APPROVED

Sohil
 9/24/2018 11:39:04 AM

Quant Time: Sep 24 07:51:46 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	114027	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	445977	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	193888	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	305320	20.00	ng	0.00
75) Chrysene-d12	13.83	240	235185	20.00	ng	-0.01
86) Perylene-d12	15.24	264	206496	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	143411	20.72	ng	-0.01
7) Phenol-d6	6.33	99	187906	21.27	ng	-0.02
23) Nitrobenzene-d5	7.26	82	111742	14.79	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	33208	17.37	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	222014	17.27	ng	-0.01
78) Terphenyl-d14	12.79	244	149921	15.64	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	52079	3.693	ng	Qvalue # 98
70) Phenanthrene	11.23	178	44931	2.947	ng	99
74) Fluoranthene	12.42	202	74694	4.585	ng	96
77) Pyrene	12.64	202	70833	4.202	ng	98
80) Benzo(a)anthracene	13.83	228	36445	2.573	ng	96
82) Chrysene	13.86	228	36175	2.576	ng	94
87) Benzo(b)fluoranthene	14.84	252	38142m	3.361	ng	
89) Benzo(a)pyrene	15.17	252	25623	2.506	ng	# 85

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

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