

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF092818\
 Data File : BF109433.D
 Acq On : 28 Sep 2018 16:46
 Operator : JU/SJ
 Sample : J4776-09 5X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-092718-C

Manual Integrations
 APPROVED

Sohil
 10/1/2018 4:19:58 PM

Quant Time: Oct 01 08:59:47 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	127334	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	461708	20.00	ng	0.00
38) Acenaphthene-d10	9.74	164	186943	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	312383	20.00	ng	0.00
75) Chrysene-d12	13.84	240	256611	20.00	ng	0.00
86) Perylene-d12	15.25	264	206876	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	113188	14.64	ng	0.00
7) Phenol-d6	6.34	99	144918	14.69	ng	-0.02
23) Nitrobenzene-d5	7.26	82	87045	11.13	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	20429	11.09	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	155417	12.54	ng	-0.01
78) Terphenyl-d14	12.80	244	119410	11.42	ng	0.00
Target Compounds						
70) Phenanthrene	11.24	178	61107	3.918	ng	98
74) Fluoranthene	12.42	202	89234	5.354	ng	97
77) Pyrene	12.65	202	86641	4.711	ng	99
80) Benzo(a)anthracene	13.83	228	38158	2.469	ng	95
82) Chrysene	13.87	228	34045	2.222	ng	97
87) Benzo(b)fluoranthene	14.85	252	34901m	3.070	ng	
89) Benzo(a)pyrene	15.19	252	23003	2.246	ng	# 80

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

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