

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109898.D
 Acq On : 15 Oct 2018 23:56
 Operator : JU/SJ
 Sample : J5198-19 2X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-02-101218-E

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:49 AM

Quant Time: Oct 16 09:09:42 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF101518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 15 18:09:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.98	152	194950	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	654515	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	268443	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	558917	20.00	ng	0.00
76) Chrysene-d12	14.15	240	546118	20.00	ng	0.00
87) Perylene-d12	15.68	264	348700	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	570644	46.40	ng	0.00
7) Phenol-d6	6.59	99	651949	42.69	ng	0.00
23) Nitrobenzene-d5	7.53	82	344394	35.45	ng	-0.01
42) 2,4,6-Tribromophenol	10.80	330	119557	51.44	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	551427	32.78	ng	0.00
79) Terphenyl-d14	13.09	244	669941	25.76	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.72	163	60361	3.176	ng	99
71) Phenanthrene	11.53	178	57872	2.000	ng	99
75) Fluoranthene	12.72	202	149155	5.160	ng	97
78) Pyrene	12.96	202	180311	4.495	ng	98
81) Benzo(a)anthracene	14.14	228	69440	2.159	ng	# 90
83) Chrysene	14.18	228	66301	2.165	ng	# 95
88) Benzo(b)fluoranthene	15.22	252	57535m	2.862	ng	
90) Benzo(a)pyrene	15.61	252	42712	2.310	ng	# 76

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

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