

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109897.D
 Acq On : 15 Oct 2018 23:28
 Operator : JU/SJ
 Sample : J5198-11 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:46 AM

Quant Time: Oct 16 02:30:31 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	240326	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	873980	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	349660	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	628791	20.00	ng	0.00
76) Chrysene-d12	14.15	240	634719	20.00	ng	0.00
87) Perylene-d12	15.68	264	434293	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	802380	52.92	ng	0.00
7) Phenol-d6	6.59	99	970838	51.57	ng	0.00
23) Nitrobenzene-d5	7.53	82	546168	42.10	ng	-0.01
42) 2,4,6-Tribromophenol	10.80	330	165399	54.64	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	796312	36.34	ng	0.00
79) Terphenyl-d14	13.09	244	726465	24.03	ng	0.00
Target Compounds						
50) Dimethylphthalate	9.72	163	95037	3.838	ng	99
71) Phenanthrene	11.53	178	74279	2.282	ng	97
75) Fluoranthene	12.72	202	162770	5.005	ng	97
78) Pyrene	12.95	202	150757	3.234	ng	98
83) Chrysene	14.17	228	77286	2.172	ng	96
88) Benzo(b)fluoranthene	15.22	252	79642m	3.181	ng	
90) Benzo(a)pyrene	15.60	252	47620	2.068	ng	# 88

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

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