

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109893.D
 Acq On : 15 Oct 2018 21:41
 Operator : JU/SJ
 Sample : J5198-17
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:41 AM

Quant Time: Oct 16 02:07:30 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	209375	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	715512	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	270854	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	510066	20.00	ng	0.00
76) Chrysene-d12	14.15	240	562655	20.00	ng	0.00
87) Perylene-d12	15.67	264	425433	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	1020795	77.28	ng	0.00
7) Phenol-d6	6.59	99	1232418	75.14	ng	0.00
23) Nitrobenzene-d5	7.53	82	693770	65.32	ng	-0.01
42) 2,4,6-Tribromophenol	10.80	330	204892	87.37	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1038842	61.20	ng	0.00
79) Terphenyl-d14	13.09	244	1061994	39.64	ng	0.00
Target Compounds						
10) Phenol	6.60	94	40290	2.275	ng	Qvalue # 76
50) Dimethylphthalate	9.72	163	100656	5.248	ng	98
71) Phenanthrene	11.53	178	138777	5.255	ng	98
75) Fluoranthene	12.72	202	380978	14.442	ng	98
78) Pyrene	12.95	202	360060	8.713	ng	98
81) Benzo(a)anthracene	14.14	228	193170	5.829	ng	98
83) Chrysene	14.17	228	162360	5.147	ng	97
86) Indeno(1,2,3-cd)pyrene	17.22	276	59565	2.345	ng	93
88) Benzo(b)fluoranthene	15.22	252	186776m	7.615	ng	
89) Benzo(k)fluoranthene	15.24	252	56056m	2.319	ng	
90) Benzo(a)pyrene	15.60	252	123707	5.484	ng	# 95
92) Benzo(a,h,i)perylene	17.69	276	58090	2.876	ng	# 89

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

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