

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101518\
 Data File : BF109892.D
 Acq On : 15 Oct 2018 21:14
 Operator : JU/SJ
 Sample : J5198-15
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/16/2018 9:57:38 AM

Quant Time: Oct 16 02:00:34 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	6.97	152	245133	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	860242	20.00	ng	0.00
39) Acenaphthene-d10	10.02	164	321726	20.00	ng	0.00
64) Phenanthrene-d10	11.51	188	532628	20.00	ng	0.00
76) Chrysene-d12	14.15	240	562539	20.00	ng	0.00
87) Perylene-d12	15.67	264	420937	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	1249643	80.81	ng	0.00
7) Phenol-d6	6.59	99	1520624	79.19	ng	0.00
23) Nitrobenzene-d5	7.54	82	894696	70.06	ng	0.00
42) 2,4,6-Tribromophenol	10.80	330	234079	84.04	ng	0.00
45) 2-Fluorobiphenyl	9.33	172	1307852	64.87	ng	0.00
79) Terphenyl-d14	13.09	244	1129006	42.15	ng	0.00
Target Compounds						
10) Phenol	6.60	94	51457	2.482	ng	Qvalue # 60
50) Dimethylphthalate	9.72	163	160396	7.041	ng	99
71) Phenanthrene	11.53	178	56578	2.052	ng	98
75) Fluoranthene	12.72	202	137703	4.999	ng	100
78) Pyrene	12.95	202	162513	3.933	ng	98
81) Benzo(a)anthracene	14.14	228	77322	2.334	ng	# 92
83) Chrysene	14.17	228	73515	2.331	ng	# 95
88) Benzo(b)fluoranthene	15.22	252	83895m	3.457	ng	
90) Benzo(a)pyrene	15.60	252	53990	2.419	ng	# 81

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

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