

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100518\
 Data File : BF109630.D
 Acq On : 6 Oct 2018 1:17
 Operator : JU/SJ
 Sample : J5263-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 GRID18-COMP-100318

Manual Integrations
 APPROVED

Sohil
 10/8/2018 12:21:44 PM

Quant Time: Oct 06 07:06:34 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	89739	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	333470	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	133603	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	219514	20.00	ng	0.00
75) Chrysene-d12	13.83	240	226819	20.00	ng	-0.01
86) Perylene-d12	15.23	264	197092	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	458288	84.11	ng	0.01
7) Phenol-d6	6.35	99	608404	87.51	ng	0.00
23) Nitrobenzene-d5	7.26	82	383237	67.84	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	114482	86.92	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	563235	63.58	ng	-0.02
78) Terphenyl-d14	12.79	244	466281	50.45	ng	-0.02
Target Compounds						
10) Phenol	6.36	94	30657	3.894	ng	Qvalue # 72
49) Dimethylphthalate	9.45	163	70561	7.261	ng	99
70) Phenanthrene	11.23	178	31352	2.861	ng	100
74) Fluoranthene	12.42	202	54979	4.694	ng	97
77) Pyrene	12.64	202	47682	2.933	ng	97
80) Benzo(a)anthracene	13.82	228	33724	2.468	ng	93
82) Chrysene	13.86	228	28983	2.140	ng	97
87) Benzo(b)fluoranthene	14.84	252	29519m	2.726	ng	
89) Benzo(a)pyrene	15.17	252	22291	2.284	ng	94

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

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