

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
 Data File : BF109848.D
 Acq On : 12 Oct 2018 21:51
 Operator : JU/SJ
 Sample : J5314-03
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 DOCUMENTATION-20

Manual Integrations
 APPROVED

Sohil
 10/15/2018 10:28:44 AM

Quant Time: Oct 15 07:05:45 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	87397	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	349759	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	142549	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	226358	20.00	ng	0.00
75) Chrysene-d12	13.83	240	191985	20.00	ng	0.00
86) Perylene-d12	15.23	264	149355	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	453758	92.16	ng	0.00
7) Phenol-d6	6.35	99	619740	94.22	ng	-0.01
23) Nitrobenzene-d5	7.26	82	405073	63.84	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	111286	71.65	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	614797	59.40	ng	-0.01
78) Terphenyl-d14	12.78	244	408409	41.18	ng	0.00
Target Compounds						
10) Phenol	6.35	94	19535	2.590	ng	Qvalue # 71
49) Dimethylphthalate	9.45	163	77560	7.419	ng	99
70) Phenanthrene	11.22	178	162983	14.388	ng	99
71) Anthracene	11.28	178	66735	5.699	ng	99
74) Fluoranthene	12.41	202	350557	29.623	ng	98
77) Pyrene	12.63	202	301784	21.455	ng	98
80) Benzo(a)anthracene	13.82	228	166040	15.672	ng	97
82) Chrysene	13.86	228	135654	12.190	ng	96
85) Indeno(1,2,3-cd)pyrene	16.58	276	51418	6.455	ng	# 93
87) Benzo(b)fluoranthene	14.84	252	126594m	15.338	ng	
88) Benzo(k)fluoranthene	14.85	252	56501m	6.813	ng	
89) Benzo(a)pyrene	15.17	252	95442	12.743	ng	98
91) Benzo(a,h,i)perylene	16.98	276	50758	8.263	ng	99

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

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