

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
 Data File : BF109702.D
 Acq On : 9 Oct 2018 18:35
 Operator : JU/SJ
 Sample : J5201-01 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-100818-A

Manual Integrations
 APPROVED

Sohil
 10/10/2018 11:37:08 AM

Quant Time: Oct 09 19:15:29 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	87772	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	383683	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	165988	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	264794	20.00	ng	0.00
75) Chrysene-d12	13.83	240	217467	20.00	ng	0.00
86) Perylene-d12	15.23	264	188389	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	118421	23.95	ng	-0.01
7) Phenol-d6	6.34	99	248615	37.64	ng	-0.02
23) Nitrobenzene-d5	7.26	82	147055	21.13	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	60459	33.43	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	323548	26.85	ng	-0.01
78) Terphenyl-d14	12.78	244	239705	21.34	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	33938	2.788	ng	99
57) Fluorene	10.27	166	29627	2.757	ng	97
70) Phenanthrene	11.23	178	250339	18.892	ng	98
71) Anthracene	11.28	178	82776	6.043	ng	98
72) Carbazole	11.45	167	41784	3.145	ng	96
74) Fluoranthene	12.41	202	414089	29.912	ng	97
77) Pyrene	12.63	202	332214	20.851	ng	98
80) Benzo(a)anthracene	13.82	228	174022	14.501	ng	97
82) Chrysene	13.85	228	157587	12.502	ng	96
85) Indeno(1,2,3-cd)pyrene	16.57	276	75435m	8.361	ng	
87) Benzo(b)fluoranthene	14.83	252	178967m	17.191	ng	
88) Benzo(k)fluoranthene	14.86	252	74593m	7.131	ng	
89) Benzo(a)pyrene	15.17	252	128654	13.618	ng	96
91) Benzo(a,h,i)perylene	16.98	276	64844	8.369	ng	99

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

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