

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101218\
 Data File : BF109847.D
 Acq On : 12 Oct 2018 21:23
 Operator : JU/SJ
 Sample : J5394-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-GF-02-027-B

Manual Integrations
 APPROVED

Sohil
 10/15/2018 10:28:41 AM

Quant Time: Oct 15 07:02:21 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	97321	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	395896	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	168600	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	259199	20.00	ng	0.00
75) Chrysene-d12	13.83	240	200156	20.00	ng	0.00
86) Perylene-d12	15.23	264	169508	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	535033	97.58	ng	0.00
7) Phenol-d6	6.35	99	731615	99.89	ng	-0.01
23) Nitrobenzene-d5	7.26	82	472205	65.75	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	150166	81.74	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	779693	63.69	ng	0.00
78) Terphenyl-d14	12.78	244	521615	50.45	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.44	163	98686	7.981	ng	99
70) Phenanthrene	11.22	178	32076	2.473	ng	98
71) Anthracene	11.28	178	27622m	2.060	ng	
74) Fluoranthene	12.40	202	86260	6.366	ng	97
77) Pyrene	12.63	202	78502	5.353	ng	98
80) Benzo(a)anthracene	13.82	228	51793	4.689	ng	97
82) Chrysene	13.86	228	51196	4.413	ng	97
85) Indeno(1,2,3-cd)pyrene	16.59	276	21633	2.605	ng	93
87) Benzo(b)fluoranthene	14.84	252	52351m	5.589	ng	
88) Benzo(k)fluoranthene	14.86	252	27022m	2.871	ng	
89) Benzo(a)pyrene	15.17	252	40885	4.810	ng	# 95
91) Benzo(a,h,i)perylene	16.99	276	17960	2.576	ng	99

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

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