

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100418\
 Data File : BF109586.D
 Acq On : 4 Oct 2018 15:04
 Operator : JU/SJ
 Sample : J5204-03 2X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-04-100318-B

Manual Integrations
 APPROVED

Sohil
 10/5/2018 11:56:29 AM

Quant Time: Oct 04 16:49:51 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	102419	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	410452	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	183177	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	294624	20.00	ng	0.00
75) Chrysene-d12	13.84	240	216587	20.00	ng	0.00
86) Perylene-d12	15.25	264	196327	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	228203	36.70	ng	0.00
7) Phenol-d6	6.34	99	326766	41.18	ng	-0.02
23) Nitrobenzene-d5	7.26	82	202707	29.15	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	64064	35.48	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	382172	31.47	ng	-0.02
78) Terphenyl-d14	12.79	244	259415	29.39	ng	-0.01
Target Compounds						
48) Acenaphthylene	9.59	152	88791	4.936	ng	99
49) Dimethylphthalate	9.45	163	41391	3.107	ng	# 98
51) Acenaphthene	9.76	154	46916	4.314	ng	99
54) Dibenzofuran	9.94	168	79634	4.924	ng	97
57) Fluorene	10.28	166	145271	12.589	ng	99
70) Phenanthrene	11.24	178	787077	53.504	ng	99
71) Anthracene	11.29	178	341402	22.882	ng	99
72) Carbazole	11.45	167	68642	4.624	ng	98
74) Fluoranthene	12.43	202	1007545	64.090	ng	97
77) Pyrene	12.66	202	903187	58.186	ng	100
80) Benzo(a)anthracene	13.84	228	523994	40.166	ng	96
82) Chrysene	13.87	228	459705	35.553	ng	98
85) Indeno(1,2,3-cd)pyrene	16.62	276	182164	17.381	ng	# 87
87) Benzo(b)fluoranthene	14.86	252	533648m	49.467	ng	
88) Benzo(k)fluoranthene	14.88	252	179627m	17.547	ng	
89) Benzo(a)pyrene	15.20	252	378691	38.956	ng	97
90) Dibenzo(a,h)anthracene	16.62	278	50963	6.390	ng	# 90
91) Benzo(g,h,i)perylene	17.02	276	163289	20.833	ng	# 89

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

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