

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF062918\
 Data File : BF106974.D
 Acq On : 29 Jun 2018 15:45
 Operator : JU/SJ
 Sample : J3677-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 147-149-B1(0-5)

Manual Integrations
 APPROVED

Sohil
 6/30/2018 10:58:04 AM

Quant Time: Jun 29 16:35:25 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF061918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 29 11:40:02 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	80961	20.00	ng	0.00
21) Naphthalene-d8	8.24	136	299654	20.00	ng	-0.01
38) Acenaphthene-d10	10.00	164	137484	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	237647	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	214554	20.00	ng	-0.02
86) Perylene-d12	15.71	264	165943	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	413718	86.53	ng	0.00
7) Phenol-d6	6.60	99	521330	87.31	ng	-0.02
23) Nitrobenzene-d5	7.52	82	316531	58.70	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	145034	91.02	ng	-0.01
44) 2-Fluorobiphenyl	9.31	172	587825	69.72	ng	-0.02
78) Terphenyl-d14	13.07	244	560482	63.28	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.70	163	83749	8.122	ng	99
51) Acenaphthene	10.02	154	18462	2.153	ng	99
70) Phenanthrene	11.51	178	175434	15.305	ng	98
71) Anthracene	11.57	178	47381	4.050	ng	95
74) Fluoranthene	12.71	202	247837	19.617	ng	95
77) Pyrene	12.94	202	216480	15.301	ng	98
80) Benzo(a)anthracene	14.14	228	127529	9.753	ng	99
82) Chrysene	14.18	228	115015	9.560	ng	96
85) Indeno(1,2,3-cd)pyrene	17.28	276	43714	4.282	ng	# 87
87) Benzo(b)fluoranthene	15.25	252	127884m	12.406	ng	
88) Benzo(k)fluoranthene	15.27	252	40514m	4.127	ng	
89) Benzo(a)pyrene	15.64	252	85454	9.325	ng	96
91) Benzo(a,h,i)perylene	17.77	276	43414	5.719	ng	# 86

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

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