

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
 Data File : BF107245.D
 Acq On : 9 Jul 2018 22:36
 Operator : JU/SJ
 Sample : J3880-13
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-17

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:47:03 PM

Quant Time: Jul 10 05:21:58 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 14:38:29 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.95	152	29072	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	96402	20.00	ng	-0.02
38) Acenaphthene-d10	10.00	164	44143	20.00	ng	-0.01
63) Phenanthrene-d10	11.49	188	91878	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	65735	20.00	ng	-0.02
86) Perylene-d12	15.69	264	52104	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	186549	108.66	ng	0.00
7) Phenol-d6	6.60	99	218831	102.07	ng	-0.02
23) Nitrobenzene-d5	7.51	82	117884	67.96	ng	-0.02
41) 2,4,6-Tribromophenol	10.80	330	56543	110.52	ng	0.00
44) 2-Fluorobiphenyl	9.30	172	224111	86.13	ng	-0.02
78) Terphenyl-d14	13.07	244	293685	108.22	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.70	163	53624	16.197	ng	Qvalue 98
70) Phenanthrene	11.51	178	38913	8.781	ng	97
71) Anthracene	11.57	178	16375	3.620	ng	97
74) Fluoranthene	12.71	202	89568	18.338	ng	96
77) Pyrene	12.95	202	82612	19.058	ng	97
80) Benzo(a)anthracene	14.14	228	38887	9.706	ng	97
82) Chrysene	14.18	228	34364	9.323	ng	98
85) Indeno(1,2,3-cd)pyrene	17.29	276	17658	5.645	ng	# 88
87) Benzo(b)fluoranthene	15.23	252	42345m	13.083	ng	
88) Benzo(k)fluoranthene	15.26	252	9802m	3.180	ng	
89) Benzo(a)pyrene	15.62	252	26162	9.092	ng	# 91
90) Dibenzo(a,h)anthracene	17.28	278	5487	2.250	ng	# 63
91) Benzo(a,h,i)perylene	17.79	276	17835	7.483	ng	# 90

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

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