

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070918\
 Data File : BF107243.D
 Acq On : 9 Jul 2018 21:41
 Operator : JU/SJ
 Sample : J3879-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SURCHARGE-SP-10

Manual Integrations
 APPROVED

Sohil
 7/10/2018 2:46:58 PM

Quant Time: Jul 10 04:56:25 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 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	29710	20.00	ng	-0.01
21) Naphthalene-d8	8.23	136	98021	20.00	ng	-0.02
38) Acenaphthene-d10	9.99	164	46868	20.00	ng	-0.02
63) Phenanthrene-d10	11.49	188	102183	20.00	ng	-0.01
75) Chrysene-d12	14.15	240	69928	20.00	ng	-0.02
86) Perylene-d12	15.69	264	53279	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.60	112	196299	111.88	ng	0.00
7) Phenol-d6	6.60	99	226927	103.57	ng	-0.02
23) Nitrobenzene-d5	7.51	82	126316	71.62	ng	-0.02
41) 2,4,6-Tribromophenol	10.79	330	64660	119.03	ng	-0.01
44) 2-Fluorobiphenyl	9.30	172	238628	86.43	ng	-0.02
78) Terphenyl-d14	13.07	244	326374	113.06	ng	-0.01
Target Compounds						
48) Acenaphthylene	9.85	152	9304	2.004	ng	99
49) Dimethylphthalate	9.70	163	48433	13.778	ng	98
70) Phenanthrene	11.51	178	65187	13.227	ng	96
71) Anthracene	11.57	178	22722	4.517	ng	98
74) Fluoranthene	12.71	202	145809	26.842	ng	95
77) Pyrene	12.95	202	138017	29.930	ng	99
80) Benzo(a)anthracene	14.14	228	62988	14.779	ng	98
82) Chrysene	14.18	228	57145	14.574	ng	97
85) Indeno(1,2,3-cd)pyrene	17.29	276	27435	8.245	ng	# 86
87) Benzo(b)fluoranthene	15.23	252	62562m	18.903	ng	
88) Benzo(k)fluoranthene	15.26	252	16242m	5.153	ng	
89) Benzo(a)pyrene	15.63	252	41645	14.154	ng	94
90) Dibenzo(a,h)anthracene	17.29	278	7865	3.154	ng	# 73
91) Benzo(g,h,i)perylene	17.79	276	27230	11.173	ng	# 88

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

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