

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF070618\
 Data File : BF107179.D
 Acq On : 6 Jul 2018 18:07
 Operator : JU/SJ
 Sample : J3820-01
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 BU-03-070318

Manual Integrations
 APPROVED

Sohil
 7/9/2018 1:58:30 PM

Quant Time: Jul 09 04:10:20 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	6.94	152	61562	20.00	ng	-0.02
21) Naphthalene-d8	8.23	136	231634	20.00	ng	-0.02
38) Acenaphthene-d10	9.98	164	102832	20.00	ng	-0.02
63) Phenanthrene-d10	11.48	188	167896	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	173817	20.00	ng	-0.03
86) Perylene-d12	15.67	264	134454	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.59	112	330395	90.88	ng	-0.01
7) Phenol-d6	6.59	99	429194	94.53	ng	-0.02
23) Nitrobenzene-d5	7.51	82	262494	62.98	ng	-0.02
41) 2,4,6-Tribromophenol	10.78	330	70085	58.80	ng	-0.02
44) 2-Fluorobiphenyl	9.30	172	462137	74.19	ng	-0.02
78) Terphenyl-d14	13.07	244	451129	62.87	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.70	163	105120	13.630	ng	# 99
57) Fluorene	10.54	166	15278	2.192	ng	# 92
70) Phenanthrene	11.51	178	109082	13.470	ng	97
71) Anthracene	11.56	178	32353	3.914	ng	97
74) Fluoranthene	12.70	202	208801	23.394	ng	94
77) Pyrene	12.93	202	194856	17.000	ng	98
80) Benzo(a)anthracene	14.12	228	91500	8.637	ng	97
82) Chrysene	14.16	228	93191	9.562	ng	98
85) Indeno(1,2,3-cd)pyrene	17.24	276	37545	4.539	ng	# 91
87) Benzo(b)fluoranthene	15.21	252	106539m	12.756	ng	
88) Benzo(k)fluoranthene	15.24	252	30357m	3.817	ng	
89) Benzo(a)pyrene	15.60	252	67428	9.081	ng	96
91) Benzo(a,h,i)perylene	17.73	276	41012	6.668	ng	# 89

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

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