

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF071118\
 Data File : BF107303.D
 Acq On : 11 Jul 2018 16:38
 Operator : JU/SJ
 Sample : J3933-02DL 10X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EP-2-4DL

Manual Integrations
 APPROVED

Sohil
 7/12/2018 2:19:26 PM

Quant Time: Jul 11 18:11:39 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	6.93	152	59708	20.00	ng	-0.03
21) Naphthalene-d8	8.21	136	228025	20.00	ng	-0.04
38) Acenaphthene-d10	9.98	164	101545	20.00	ng	-0.03
63) Phenanthrene-d10	11.48	188	173287	20.00	ng	-0.02
75) Chrysene-d12	14.14	240	187942	20.00	ng	-0.02
86) Perylene-d12	15.68	264	155518	20.00	ng	-0.04
System Monitoring Compounds						
5) 2-Fluorophenol	5.57	112	33267	9.43	ng	-0.03
7) Phenol-d6	6.58	99	42424	9.63	ng	-0.04
23) Nitrobenzene-d5	7.50	82	25004	6.09	ng	-0.04
41) 2,4,6-Tribromophenol	10.78	330	8818	7.49	ng	-0.02
44) 2-Fluorobiphenyl	0.00	172	0d	0.00	ng	
78) Terphenyl-d14	13.06	244	50643	6.53	ng	-0.02
Target Compounds						
51) Acenaphthene	10.01	154	15477	2.444	ng	Qvalue 98
57) Fluorene	10.53	166	16883	2.453	ng	# 98
70) Phenanthrene	11.50	178	364217	43.577	ng	99
71) Anthracene	11.55	178	55897	6.552	ng	99
74) Fluoranthene	12.70	202	388913	42.217	ng	96
77) Pyrene	12.94	202	527115	42.532	ng	99
80) Benzo(a)anthracene	14.13	228	320489	27.979	ng	98
82) Chrysene	14.17	228	315294	29.920	ng	98
85) Indeno(1,2,3-cd)pyrene	17.28	276	86521	9.675	ng	# 87
87) Benzo(b)fluoranthene	15.22	252	294198m	30.453	ng	
88) Benzo(k)fluoranthene	15.25	252	62582m	6.803	ng	
89) Benzo(a)pyrene	15.62	252	200173m	23.308	ng	
90) Dibenzo(a,h)anthracene	17.28	278	30756m	4.226	ng	
91) Benzo(a,h,i)perylene	17.77	276	104954	14.753	ng	# 88

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

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