

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111303.D
 Acq On : 12 Dec 2018 3:49
 Operator : JU/SJ
 Sample : J6193-05 5X
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-121018-C

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:30:22 PM

Quant Time: Dec 12 04:51:22 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 10 01:15:27 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	250894	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	978324	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	422766	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	665318	20.00	ng	0.00
76) Chrysene-d12	13.95	240	588128	20.00	ng	0.00
87) Perylene-d12	15.39	264	426907	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	393464	24.76	ng	0.00
7) Phenol-d6	6.42	99	498657	25.25	ng	-0.01
23) Nitrobenzene-d5	7.35	82	287870	16.50	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	80378	21.46	ng	-0.01
45) 2-Fluorobiphenyl	9.15	172	515065	14.25	ng	-0.01
79) Terphenyl-d14	12.90	244	408876	16.68	ng	0.00
Target Compounds						Qvalue
50) Dimethylphthalate	9.55	163	89166	2.858	ng	98
71) Phenanthrene	11.34	178	91570	2.729	ng	94
75) Fluoranthene	12.53	202	223200	7.281	ng	97
78) Pyrene	12.75	202	201389	4.792	ng	97
81) Benzo(a)anthracene	13.94	228	129448	3.945	ng	99
83) Chrysene	13.97	228	112139	3.315	ng	99
88) Benzo(b)fluoranthene	14.97	252	117078m	4.959	ng	
90) Benzo(a)pyrene	15.33	252	73565	3.384	ng	97

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

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