

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111610.D
 Acq On : 20 Dec 2018 00:29
 Operator : JU/SJ
 Sample : J6196-09 2X
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-01-121818-B

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:22 PM

Quant Time: Dec 20 04:29:21 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 19 17:39:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	235028	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	878268	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	423229	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	679255	20.00	ng	0.00
76) Chrysene-d12	14.04	240	551962	20.00	ng	0.00
87) Perylene-d12	15.52	264	418174	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	453179	32.87	ng	-0.01
7) Phenol-d6	6.51	99	609172	34.71	ng	-0.01
23) Nitrobenzene-d5	7.45	82	344487	22.83	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	148339	29.66	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	689362	23.74	ng	0.00
79) Terphenyl-d14	12.99	244	587331	20.18	ng	0.00
Target Compounds						
10) Phenol	6.52	94	47329	2.390	ng	95
50) Dimethylphthalate	9.64	163	101364	3.275	ng	# 98
71) Phenanthrene	11.43	178	224461	6.231	ng	94
72) Anthracene	11.49	178	72338	2.001	ng	95
75) Fluoranthene	12.62	202	225068	6.170	ng	95
78) Pyrene	12.85	202	199227	5.111	ng	97
81) Benzo(a)anthracene	14.03	228	104439	3.315	ng	98
83) Chrysene	14.07	228	84307	2.723	ng	96
88) Benzo(b)fluoranthene	15.09	252	97704m	3.998	ng	
90) Benzo(a)pyrene	15.45	252	65815	2.964	ng	98

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

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