

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121018\
 Data File : BF111263.D
 Acq On : 11 Dec 2018 00:54
 Operator : JU/SJ
 Sample : J6196-01 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-03-120718-A

Manual Integrations
 APPROVED

Sohil
 12/11/2018 2:23:47 PM

Quant Time: Dec 11 06:23:36 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	464366	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	1690714	20.00	ng	0.00
39) Acenaphthene-d10	9.84	164	648726	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1011554	20.00	ng	0.00
76) Chrysene-d12	13.96	240	751429	20.00	ng	0.00
87) Perylene-d12	15.40	264	563585	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1551967	52.77	ng	0.00
7) Phenol-d6	6.43	99	1889409	51.70	ng	0.00
23) Nitrobenzene-d5	7.36	82	1090012	36.16	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	254542	44.30	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1589753	34.68	ng	0.00
79) Terphenyl-d14	12.90	244	1211486	38.67	ng	0.00
Target Compounds						
10) Phenol	6.44	94	106421	2.490	ng	93
50) Dimethylphthalate	9.55	163	170881	3.570	ng	99
71) Phenanthrene	11.35	178	111442	2.184	ng	93
75) Fluoranthene	12.53	202	213819	4.588	ng	99
78) Pyrene	12.76	202	195716	3.645	ng	97
81) Benzo(a)anthracene	13.95	228	94373	2.251	ng	# 91
88) Benzo(b)fluoranthene	14.99	252	90509m	2.904	ng	

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

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