

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121118\
 Data File : BF111279.D
 Acq On : 11 Dec 2018 15:26
 Operator : JU/SJ
 Sample : J6197-01 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 PL-02-120718

Manual Integrations
 APPROVED

Sohil
 12/12/2018 5:30:07 PM

Quant Time: Dec 11 17:39:48 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	341114	20.00	ng	0.00
21) Naphthalene-d8	8.08	136	1244021	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	538313	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	834029	20.00	ng	0.00
76) Chrysene-d12	13.96	240	445000	20.00	ng	0.00
87) Perylene-d12	15.40	264	348311	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.41	112	863178	39.96	ng	0.00
7) Phenol-d6	6.43	99	1456409	54.25	ng	0.00
23) Nitrobenzene-d5	7.35	82	827412	37.31	ng	-0.01
42) 2,4,6-Tribromophenol	10.62	330	175832	36.87	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1168934	29.84	ng	0.00
79) Terphenyl-d14	12.90	244	870651	46.93	ng	0.00
Target Compounds						
10) Phenol	6.44	94	69961	2.229	ng	93
50) Dimethylphthalate	9.55	163	86464	2.177	ng	98
71) Phenanthrene	11.34	178	139367	3.313	ng	92
75) Fluoranthene	12.53	202	161259	4.196	ng	98
78) Pyrene	12.76	202	159123	5.004	ng	95
81) Benzo(a)anthracene	13.94	228	57236	2.305	ng	# 90
88) Benzo(b)fluoranthene	14.98	252	49159m	2.552	ng	
90) Benzo(a)pyrene	15.33	252	37703	2.126	ng	# 79

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

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