

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121918\
 Data File : BF111613.D
 Acq On : 20 Dec 2018 1:52
 Operator : JU/SJ
 Sample : J6194-07 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OK-01-121818-A

Manual Integrations
 APPROVED

Sohil
 12/20/2018 12:42:33 PM

Quant Time: Dec 20 04:37:37 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	6.90	152	189894	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	705477	20.00	ng	0.00
39) Acenaphthene-d10	9.93	164	317384	20.00	ng	0.00
64) Phenanthrene-d10	11.41	188	542519	20.00	ng	0.00
76) Chrysene-d12	14.04	240	496637	20.00	ng	0.00
87) Perylene-d12	15.52	264	344550	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	203188	18.24	ng	-0.01
7) Phenol-d6	6.51	99	256515	18.09	ng	-0.01
23) Nitrobenzene-d5	7.45	82	134228	11.07	ng	-0.01
42) 2,4,6-Tribromophenol	10.71	330	55021	14.67	ng	-0.01
45) 2-Fluorobiphenyl	9.25	172	247040	11.35	ng	0.00
79) Terphenyl-d14	12.99	244	237390	9.06	ng	0.00
Target Compounds						
71) Phenanthrene	11.43	178	62982	2.189	ng	93
75) Fluoranthene	12.62	202	125432	4.305	ng	95
78) Pyrene	12.85	202	118476	3.378	ng	96
81) Benzo(a)anthracene	14.03	228	61940	2.185	ng	93
83) Chrysene	14.07	228	55732	2.001	ng	97
88) Benzo(b)fluoranthene	15.09	252	63181m	3.138	ng	
90) Benzo(a)pyrene	15.46	252	45049	2.462	ng	# 92

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

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