

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF121418\
 Data File : BF111419.D
 Acq On : 14 Dec 2018 15:27
 Operator : JU/SJ
 Sample : J6319-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 059_BT-3-1

Manual Integrations
 APPROVED

Sohil
 12/17/2018 2:36:22 PM

Quant Time: Dec 14 16:43:17 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF121418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 14 13:26:17 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.79	152	170201	20.00	ng	0.00
21) Naphthalene-d8	8.07	136	683392	20.00	ng	0.00
39) Acenaphthene-d10	9.83	164	344100	20.00	ng	0.00
64) Phenanthrene-d10	11.31	188	612752	20.00	ng	0.00
76) Chrysene-d12	13.94	240	385891	20.00	ng	0.00
87) Perylene-d12	15.39	264	342382	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1412789	138.13	ng	0.00
7) Phenol-d6	6.44	99	1809202	132.98	ng	0.00
23) Nitrobenzene-d5	7.35	82	1109216	96.65	ng	0.00
42) 2,4,6-Tribromophenol	10.62	330	402049	130.43	ng	0.00
45) 2-Fluorobiphenyl	9.15	172	1756140	78.83	ng	0.00
79) Terphenyl-d14	12.90	244	1532599	93.26	ng	0.00
Target Compounds						
10) Phenol	6.45	94	122760	8.100	ng	88
50) Dimethylphthalate	9.54	163	167843	6.773	ng	98
75) Fluoranthene	12.52	202	269711	8.732	ng	98
78) Pyrene	12.75	202	370935	13.634	ng	95
81) Benzo(a)anthracene	13.93	228	116901	5.571	ng	99
83) Chrysene	13.97	228	109473	4.953	ng	95
86) Indeno(1,2,3-cd)pyrene	16.79	276	61770	3.872	ng	# 91
88) Benzo(b)fluoranthene	14.97	252	120152m	6.019	ng	
89) Benzo(k)fluoranthene	14.99	252	46024m	2.413	ng	
90) Benzo(a)pyrene	15.32	252	101793	5.655	ng	98
92) Benzo(g,h,i)perylene	17.21	276	60834	4.365	ng	94

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

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