

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF120618\
 Data File : BF111165.D
 Acq On : 7 Dec 2018 3:38
 Operator : JU/SJ
 Sample : J6199-01 2X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SU-01-120518-A

Manual Integrations
 APPROVED

mohammad
 12/10/2018 12:19:27 PM

Quant Time: Dec 07 14:11:48 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF120518.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 05 17:30:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	519573	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	1903756	20.00	ng	0.00
39) Acenaphthene-d10	9.85	164	742963	20.00	ng	0.00
64) Phenanthrene-d10	11.32	188	1141582	20.00	ng	0.00
76) Chrysene-d12	13.96	240	845521	20.00	ng	0.01
87) Perylene-d12	15.40	264	608255	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.42	112	1462452	43.92	ng	0.00
7) Phenol-d6	6.43	99	1808929	42.17	ng	0.00
23) Nitrobenzene-d5	7.36	82	1058235	32.91	ng	0.00
42) 2,4,6-Tribromophenol	10.63	330	251297	42.55	ng	0.00
45) 2-Fluorobiphenyl	9.16	172	1572708	29.86	ng	0.00
79) Terphenyl-d14	12.91	244	1175007	28.46	ng	0.00
Target Compounds						
10) Phenol	6.44	94	115956	2.479	ng	97
50) Dimethylphthalate	9.56	163	134437	2.600	ng	# 95
71) Phenanthrene	11.35	178	168495	3.068	ng	95
75) Fluoranthene	12.53	202	287852	5.357	ng	98
78) Pyrene	12.76	202	306409	4.652	ng	97
81) Benzo(a)anthracene	13.95	228	122049	2.464	ng	94
83) Chrysene	13.99	228	126359	2.569	ng	94
88) Benzo(b)fluoranthene	14.99	252	115558m	3.353	ng	
90) Benzo(a)pyrene	15.34	252	81372	2.574	ng	# 69

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

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