

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF102418\
 Data File : BF110121.D
 Acq On : 24 Oct 2018 18:05
 Operator : JU/SJ
 Sample : J5198-21 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 CL-01-102318-A

Manual Integrations
 APPROVED

Sohil
 10/25/2018 12:41:46 PM

Quant Time: Oct 25 10:04:50 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF102318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 23 15:06:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	155586	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	587025	20.00	ng	-0.02
39) Acenaphthene-d10	10.01	164	232633	20.00	ng	-0.02
64) Phenanthrene-d10	11.50	188	399137	20.00	ng	-0.01
76) Chrysene-d12	14.15	240	346394	20.00	ng	-0.01
87) Perylene-d12	15.69	264	253012	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.58	112	463120	48.47	ng	-0.02
7) Phenol-d6	6.58	99	559671	45.80	ng	-0.02
23) Nitrobenzene-d5	7.53	82	309687	31.29	ng	-0.02
42) 2,4,6-Tribromophenol	10.80	330	88801	38.54	ng	-0.02
45) 2-Fluorobiphenyl	9.32	172	533786	36.03	ng	-0.02
79) Terphenyl-d14	13.09	244	486732	29.22	ng	-0.02
Target Compounds						
10) Phenol	6.59	94	29009	2.221	ng	Qvalue # 67
50) Dimethylphthalate	9.72	163	56524	3.292	ng	99
71) Phenanthrene	11.53	178	63468	3.039	ng	98
75) Fluoranthene	12.72	202	115199	5.435	ng	99
78) Pyrene	12.95	202	119147	4.798	ng	97
81) Benzo(a)anthracene	14.14	228	55453	2.700	ng	92
83) Chrysene	14.18	228	54401	2.788	ng	97
88) Benzo(b)fluoranthene	15.24	252	63433m	4.342	ng	
90) Benzo(a)pyrene	15.62	252	41381	3.078	ng	# 85
92) Benzo(g,h,i)perylene	17.73	276	23102	2.021	ng	# 85

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

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