

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100218\
 Data File : BF109501.D
 Acq On : 2 Oct 2018 15:11
 Operator : JU/SJ
 Sample : J5200-01 2X
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-02-100118-A

Manual Integrations
 APPROVED

Sohil
 10/3/2018 9:31:01 AM

Quant Time: Oct 03 09:04:01 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF092218.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 22 13:32:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	112002	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	444553	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	211074	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	387766	20.00	ng	0.00
75) Chrysene-d12	13.84	240	279840	20.00	ng	0.00
86) Perylene-d12	15.24	264	295177	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	248693	36.57	ng	0.00
7) Phenol-d6	6.35	99	350621	40.41	ng	-0.01
23) Nitrobenzene-d5	7.26	82	219574	29.16	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	85675	41.18	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	438315	31.32	ng	-0.02
78) Terphenyl-d14	12.79	244	374736	32.86	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	48996	3.191	ng	Qvalue # 99
57) Fluorene	10.28	166	44140	3.320	ng	96
70) Phenanthrene	11.23	178	247728	12.795	ng	99
71) Anthracene	11.29	178	59161	3.013	ng	98
74) Fluoranthene	12.42	202	269693	13.035	ng	95
77) Pyrene	12.64	202	286366	14.279	ng	100
80) Benzo(a)anthracene	13.83	228	126112	7.482	ng	93
82) Chrysene	13.86	228	117279	7.020	ng	97
85) Indeno(1,2,3-cd)pyrene	16.59	276	51590	3.810	ng	# 90
87) Benzo(b)fluoranthene	14.84	252	129607m	7.991	ng	
88) Benzo(k)fluoranthene	14.87	252	43262m	2.811	ng	
89) Benzo(a)pyrene	15.18	252	103228	7.063	ng	98
91) Benzo(a,h,i)perylene	16.99	276	45774	3.884	ng	# 92

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

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