

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100218\
 Data File : BF109502.D
 Acq On : 2 Oct 2018 15:37
 Operator : JU/SJ
 Sample : J5200-03 2X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/3/2018 9:31:04 AM

Quant Time: Oct 03 04:12:29 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	109302	20.00	ng	0.00
21) Naphthalene-d8	7.99	136	426778	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	193977	20.00	ng	0.00
63) Phenanthrene-d10	11.21	188	312625	20.00	ng	0.00
75) Chrysene-d12	13.84	240	253812	20.00	ng	0.00
86) Perylene-d12	15.24	264	251620	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	194159	29.26	ng	0.00
7) Phenol-d6	6.35	99	270036	31.89	ng	-0.01
23) Nitrobenzene-d5	7.26	82	168359	23.29	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	57513	30.08	ng	-0.02
44) 2-Fluorobiphenyl	9.06	172	335413	26.08	ng	-0.02
78) Terphenyl-d14	12.79	244	244592	23.65	ng	-0.01
Target Compounds						Qvalue
49) Dimethylphthalate	9.45	163	29469	2.089	ng	100
70) Phenanthrene	11.23	178	77773	4.982	ng	99
74) Fluoranthene	12.42	202	101746	6.099	ng	98
77) Pyrene	12.64	202	114548	6.297	ng	97
80) Benzo(a)anthracene	13.83	228	55441m	3.626	ng	
82) Chrysene	13.86	228	52318	3.453	ng	96
87) Benzo(b)fluoranthene	14.84	252	65211m	4.716	ng	
89) Benzo(a)pyrene	15.18	252	47167	3.786	ng	95
91) Benzo(g,h,i)perylene	17.00	276	25674	2.556	ng	# 90

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

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