

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100918\
 Data File : BF109703.D
 Acq On : 9 Oct 2018 19:03
 Operator : JU/SJ
 Sample : J5201-03 2X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 EO-02-100818-B

Manual Integrations
 APPROVED

Sohil
 10/10/2018 11:37:11 AM

Quant Time: Oct 10 09:37:11 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF100818.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 08 16:02:11 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.70	152	74541	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	308951	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	132698	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	218532	20.00	ng	0.00
75) Chrysene-d12	13.83	240	210212	20.00	ng	0.00
86) Perylene-d12	15.23	264	206674	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	174165	41.47	ng	-0.01
7) Phenol-d6	6.34	99	280685	50.03	ng	-0.02
23) Nitrobenzene-d5	7.26	82	176340	31.46	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	53787	37.20	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	284118	29.49	ng	0.00
78) Terphenyl-d14	12.78	244	203807	18.77	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	33370	3.429	ng	100
70) Phenanthrene	11.23	178	44596	4.078	ng	99
74) Fluoranthene	12.41	202	90256	7.900	ng	99
77) Pyrene	12.63	202	80206	5.208	ng	97
80) Benzo(a)anthracene	13.82	228	44237	3.813	ng	97
82) Chrysene	13.85	228	40570	3.330	ng	98
85) Indeno(1,2,3-cd)pyrene	16.57	276	19959	2.288	ng	# 93
87) Benzo(b)fluoranthene	14.83	252	48975m	4.288	ng	
89) Benzo(a)pyrene	15.17	252	38372	3.702	ng	97
91) Benzo(g,h,i)perylene	16.98	276	20070	2.361	ng	94

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

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