

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF101018\
 Data File : BF109757.D
 Acq On : 10 Oct 2018 22:09
 Operator : JU/SJ
 Sample : J5200-07 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 OR-03-100818-B

Manual Integrations
 APPROVED

Sohil
 10/11/2018 2:02:56 PM

Quant Time: Oct 11 10:24:49 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	6.69	152	82987	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	358895	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	174238	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	312222	20.00	ng	0.00
75) Chrysene-d12	13.83	240	227458	20.00	ng	0.00
86) Perylene-d12	15.23	264	217955	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.31	112	170034	36.37	ng	-0.01
7) Phenol-d6	6.34	99	265731	42.55	ng	-0.02
23) Nitrobenzene-d5	7.26	82	168998	25.96	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	69812	36.77	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	346151	27.36	ng	-0.01
78) Terphenyl-d14	12.78	244	280292	23.85	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	45376	3.551	ng	99
70) Phenanthrene	11.22	178	109105	6.983	ng	98
74) Fluoranthene	12.41	202	110028	6.741	ng	99
77) Pyrene	12.63	202	113952	6.838	ng	98
80) Benzo(a)anthracene	13.82	228	40730	3.245	ng	95
82) Chrysene	13.85	228	46722	3.544	ng	96
87) Benzo(b)fluoranthene	14.83	252	41886m	3.478	ng	
89) Benzo(a)pyrene	15.17	252	34748	3.179	ng	# 91
91) Benzo(g,h,i)perylene	16.98	276	21569	2.406	ng	92

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

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