

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100818\
 Data File : BF109683.D
 Acq On : 9 Oct 2018 8:24
 Operator : JU/SJ
 Sample : J5309-08
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-16

Manual Integrations
 APPROVED

Sohil
 10/9/2018 2:58:05 PM

Quant Time: Oct 09 11:03:40 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	103740	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	394008	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	156554	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	264238	20.00	ng	0.00
75) Chrysene-d12	13.83	240	226002	20.00	ng	0.00
86) Perylene-d12	15.23	264	179864	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	580927	99.40	ng	0.00
7) Phenol-d6	6.35	99	801846	102.70	ng	0.00
23) Nitrobenzene-d5	7.26	82	514458	71.97	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	151680	88.92	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	868280	76.39	ng	0.00
78) Terphenyl-d14	12.78	244	686475	58.80	ng	0.00
Target Compounds						Qvalue
49) Dimethylphthalate	9.45	163	96080	8.368	ng	99
70) Phenanthrene	11.23	178	63121	4.773	ng	99
74) Fluoranthene	12.41	202	90867	6.578	ng	98
77) Pyrene	12.63	202	81661	4.932	ng	98
80) Benzo(a)anthracene	13.82	228	37210	2.984	ng	95
82) Chrysene	13.85	228	34226	2.613	ng	98
87) Benzo(b)fluoranthene	14.83	252	29072m	2.925	ng	
89) Benzo(a)pyrene	15.17	252	25737	2.853	ng	# 93

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

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