

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100818\
 Data File : BF109687.D
 Acq On : 9 Oct 2018 10:14
 Operator : JU/SJ
 Sample : J5309-17
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-25

Manual Integrations
 APPROVED

Sohil
 10/9/2018 2:58:12 PM

Quant Time: Oct 09 11:21:29 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	94378	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	350157	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	138458	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	238076	20.00	ng	0.00
75) Chrysene-d12	13.83	240	206320	20.00	ng	0.00
86) Perylene-d12	15.24	264	160498	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	565503	106.36	ng	0.00
7) Phenol-d6	6.35	99	736412	103.68	ng	0.00
23) Nitrobenzene-d5	7.26	82	471605	74.24	ng	-0.01
41) 2,4,6-Tribromophenol	10.51	330	136741	90.64	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	761467	75.74	ng	0.00
78) Terphenyl-d14	12.78	244	609530	57.19	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	144253	14.206	ng	99
74) Fluoranthene	12.41	202	45070	3.621	ng	97
77) Pyrene	12.63	202	43022	2.846	ng	98
83) Bis(2-ethylhexyl)phthalate	13.82	149	25856	3.234	ng	95
87) Benzo(b)fluoranthene	14.85	252	20493m	2.311	ng	

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

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