

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF100818\
 Data File : BF109684.D
 Acq On : 9 Oct 2018 8:51
 Operator : JU/SJ
 Sample : J5309-11
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S-19

Manual Integrations
 APPROVED

Sohil
 10/9/2018 2:58:07 PM

Quant Time: Oct 09 11:10:34 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	99328	20.00	ng	0.00
21) Naphthalene-d8	7.97	136	371347	20.00	ng	0.00
38) Acenaphthene-d10	9.72	164	155422	20.00	ng	0.00
63) Phenanthrene-d10	11.20	188	264998	20.00	ng	0.00
75) Chrysene-d12	13.83	240	215269	20.00	ng	0.00
86) Perylene-d12	15.23	264	179123	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	703488	125.72	ng	0.00
7) Phenol-d6	6.35	99	941873	126.00	ng	0.00
23) Nitrobenzene-d5	7.26	82	617246	91.63	ng	0.00
41) 2,4,6-Tribromophenol	10.51	330	195086	115.20	ng	0.00
44) 2-Fluorobiphenyl	9.05	172	1042172	92.35	ng	0.00
78) Terphenyl-d14	12.78	244	833856	74.99	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.45	163	111017	9.739	ng	99
70) Phenanthrene	11.23	178	40314	3.040	ng	99
74) Fluoranthene	12.41	202	86192	6.221	ng	98
77) Pyrene	12.63	202	82173	5.210	ng	98
80) Benzo(a)anthracene	13.82	228	33405	2.812	ng	98
82) Chrysene	13.85	228	35577m	2.851	ng	
87) Benzo(b)fluoranthene	14.83	252	36716m	3.709	ng	
89) Benzo(a)pyrene	15.17	252	25596	2.849	ng	95
91) Benzo(g,h,i)perylene	16.98	276	15602	2.118	ng	# 86

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

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