

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091818\
 Data File : BF109110.D
 Acq On : 18 Sep 2018 21:50
 Operator : JU/SJ
 Sample : J4962-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 291-11984

Manual Integrations
 APPROVED

Sohil
 9/19/2018 4:24:30 PM

Quant Time: Sep 19 05:21:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF091418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 14 13:26:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.72	152	127237	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	452929	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	187959	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	315517	20.00	ng	0.00
75) Chrysene-d12	13.85	240	326188	20.00	ng	0.00
86) Perylene-d12	15.26	264	300205	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.33	112	291546	38.06	ng	0.00
7) Phenol-d6	6.35	99	471878	47.77	ng	0.00
23) Nitrobenzene-d5	7.28	82	312121	38.65	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	29103	14.27	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	578677	48.24	ng	0.00
78) Terphenyl-d14	12.81	244	503164	36.56	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	76234	5.675	ng	99
70) Phenanthrene	11.25	178	61245	3.960	ng	98
74) Fluoranthene	12.44	202	84828	5.195	ng	94
77) Pyrene	12.66	202	94546	3.902	ng	97
80) Benzo(a)anthracene	13.85	228	43267	2.191	ng	97
82) Chrysene	13.88	228	44883	2.272	ng	95
87) Benzo(b)fluoranthene	14.86	252	43943m	2.646	ng	
89) Benzo(a)pyrene	15.20	252	32918	2.232	ng	# 95

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

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