

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF091718\
 Data File : BF109058.D
 Acq On : 17 Sep 2018 19:46
 Operator : JU/SJ
 Sample : J4946-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 RB12983RT

Manual Integrations
 APPROVED

Sohil
 9/18/2018 6:00:15 PM

Quant Time: Sep 18 04:18:47 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	108129	20.00	ng	0.00
21) Naphthalene-d8	8.00	136	357727	20.00	ng	0.00
38) Acenaphthene-d10	9.75	164	160332	20.00	ng	0.00
63) Phenanthrene-d10	11.23	188	330607	20.00	ng	0.00
75) Chrysene-d12	13.85	240	362029	20.00	ng	0.00
86) Perylene-d12	15.25	264	265430	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.32	112	401595	61.69	ng	0.00
7) Phenol-d6	6.35	99	473513	56.41	ng	0.00
23) Nitrobenzene-d5	7.28	82	268664	42.12	ng	-0.01
41) 2,4,6-Tribromophenol	10.54	330	110539	63.53	ng	0.00
44) 2-Fluorobiphenyl	9.08	172	366068	35.77	ng	0.00
78) Terphenyl-d14	12.81	244	424279	27.78	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.47	163	60462	5.276	ng	Qvalue # 98
70) Phenanthrene	11.25	178	60095	3.708	ng	96
74) Fluoranthene	12.44	202	95251	5.567	ng	96
77) Pyrene	12.66	202	94914	3.530	ng	98
80) Benzo(a)anthracene	13.84	228	45419	2.072	ng	96
82) Chrysene	13.88	228	47759	2.178	ng	95
87) Benzo(b)fluoranthene	14.86	252	48407m	3.297	ng	
89) Benzo(a)pyrene	15.19	252	30150	2.312	ng	# 95

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

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