

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092817\
 Data File : BF098928.D
 Acq On : 29 Sep 2017 1:16
 Operator : SJ/JU
 Sample : I5473-01 2X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TR-04-092717

Manual Integrations
 APPROVED

Sohil
 9/29/2017 6:04:48 PM

Quant Time: Sep 29 03:41:14 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 23 13:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.92	152	146714	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	535229	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	191239	20.00	ng	0.00
63) Phenanthrene-d10	11.44	188	288630	20.00	ng	0.00
75) Chrysene-d12	14.08	240	284348	20.00	ng	0.00
86) Perylene-d12	15.58	264	229763	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	452892	49.14	ng	0.00
7) Phenol-d6	6.53	99	524315	46.12	ng	-0.01
23) Nitrobenzene-d5	7.47	82	302927	36.30	ng	-0.01
41) 2,4,6-Tribromophenol	10.74	330	62825	40.15	ng	-0.01
44) 2-Fluorobiphenyl	9.27	172	498040	43.48	ng	-0.01
78) Terphenyl-d14	13.03	244	357735	28.61	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.66	163	83041	5.753	ng	100
70) Phenanthrene	11.46	178	33617	2.176	ng	100
74) Fluoranthene	12.66	202	60146	3.845	ng	98
77) Pyrene	12.89	202	73505	3.390	ng	98
80) Benzo(a)anthracene	14.07	228	43241	2.511	ng	# 94
82) Chrysene	14.11	228	44344m	2.705	ng	
87) Benzo(b)fluoranthene	15.13	252	51552m	3.649	ng	
89) Benzo(a)pyrene	15.51	252	36338	2.802	ng	# 89

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

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