

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092617\
 Data File : BF098837.D
 Acq On : 26 Sep 2017 17:31
 Operator : SJ/JU
 Sample : I5363-15 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-3-H-COMP

Manual Integrations
 APPROVED

Sohil
 9/27/2017 12:44:57 PM

Quant Time: Sep 27 01:58:15 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	152043	20.00	ng	0.00
21) Naphthalene-d8	8.20	136	631686	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	280207	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	452773	20.00	ng	0.00
75) Chrysene-d12	14.07	240	278502	20.00	ng	-0.01
86) Perylene-d12	15.57	264	272601	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	122498	12.83	ng	-0.02
7) Phenol-d6	6.53	99	153726	13.05	ng	-0.02
23) Nitrobenzene-d5	7.47	82	89263	9.06	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	26338	11.49	ng	-0.02
44) 2-Fluorobiphenyl	9.27	172	180193	10.74	ng	-0.01
78) Terphenyl-d14	13.02	244	118993	9.72	ng	-0.02
Target Compounds						
70) Phenanthrene	11.46	178	65539	2.704	ng	98
74) Fluoranthene	12.65	202	94870	3.866	ng	98
77) Pyrene	12.88	202	81364	3.831	ng	98
80) Benzo(a)anthracene	14.07	228	43040	2.552	ng	93
82) Chrysene	14.10	228	45824	2.854	ng	98
87) Benzo(b)fluoranthene	15.13	252	62360m	3.721	ng	
89) Benzo(a)pyrene	15.50	252	33488	2.177	ng	# 86

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

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