

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092617\
 Data File : BF098834.D
 Acq On : 26 Sep 2017 16:08
 Operator : SJ/JU
 Sample : I5363-05 10X
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 9/27/2017 12:44:49 PM

Quant Time: Sep 27 01:47:30 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	6.91	152	153485	20.00	ng	-0.01
21) Naphthalene-d8	8.20	136	652235	20.00	ng	0.00
38) Acenaphthene-d10	9.95	164	294752	20.00	ng	-0.01
63) Phenanthrene-d10	11.44	188	504760	20.00	ng	0.00
75) Chrysene-d12	14.07	240	334779	20.00	ng	-0.01
86) Perylene-d12	15.57	264	318049	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.52	112	146106	15.15	ng	-0.02
7) Phenol-d6	6.53	99	182576	15.35	ng	-0.02
23) Nitrobenzene-d5	7.47	82	104195	10.24	ng	-0.02
41) 2,4,6-Tribromophenol	10.73	330	34480	14.30	ng	-0.02
44) 2-Fluorobiphenyl	9.27	172	221947	12.57	ng	-0.01
78) Terphenyl-d14	13.02	244	172221	11.70	ng	-0.02
Target Compounds						
70) Phenanthrene	11.46	178	70612	2.613	ng	Qvalue 97
74) Fluoranthene	12.65	202	99857	3.650	ng	98
77) Pyrene	12.88	202	89678	3.513	ng	96
80) Benzo(a)anthracene	14.07	228	45619	2.250	ng	94
82) Chrysene	14.10	228	49634	2.572	ng	97
87) Benzo(b)fluoranthene	15.13	252	67599m	3.457	ng	
89) Benzo(a)pyrene	15.50	252	36246	2.019	ng	# 86

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

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