

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092217\
 Data File : BF098698.D
 Acq On : 22 Sep 2017 14:24
 Operator : SJ/JU
 Sample : I5293-07 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SP-A-COMP

Manual Integrations
 APPROVED

Sohil
 9/25/2017 6:21:08 PM

Quant Time: Sep 22 15:08:52 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 11:54:28 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.93	152	151639	20.00	ng	0.00
21) Naphthalene-d8	8.21	136	575235	20.00	ng	0.00
38) Acenaphthene-d10	9.96	164	218305	20.00	ng	0.00
63) Phenanthrene-d10	11.45	188	317847	20.00	ng	0.00
75) Chrysene-d12	14.09	240	331600	20.00	ng	0.00
86) Perylene-d12	15.58	264	326196	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.53	112	150524	15.07	ng	0.00
7) Phenol-d6	6.53	99	182005	15.05	ng	-0.01
23) Nitrobenzene-d5	7.48	82	92847	10.95	ng	-0.01
41) 2,4,6-Tribromophenol	10.75	330	22396	11.22	ng	0.00
44) 2-Fluorobiphenyl	9.28	172	182429	13.30	ng	0.00
78) Terphenyl-d14	13.03	244	127024	8.23	ng	0.00
Target Compounds						
70) Phenanthrene	11.47	178	40404	2.343	ng	98
74) Fluoranthene	12.66	202	87791	5.109	ng	98
77) Pyrene	12.89	202	90798	3.688	ng	99
80) Benzo(a)anthracene	14.07	228	40907	2.048	ng	95
82) Chrysene	14.11	228	39221	2.033	ng	97
87) Benzo(b)fluoranthene	15.14	252	47018m	2.485	ng	

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

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