

Data Path : Z:\HPCHEM1\BNA F\DATA\BF101617\
 Data File : BF099567.D
 Acq On : 17 Oct 2017 1:57
 Operator : SJ/JU
 Sample : I5752-08RE 2X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-114-3(0-3)RE

Manual Integrations
 APPROVED

Sohil
 10/17/2017 5:02:06 PM

Quant Time: Oct 17 03:08:17 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 16:25:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.84	152	101842	20.00	ng	-0.01
21) Naphthalene-d8	8.12	136	406446	20.00	ng	-0.01
38) Acenaphthene-d10	9.88	164	174318	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	272231	20.00	ng	0.00
75) Chrysene-d12	14.01	240	188423	20.00	ng	0.00
86) Perylene-d12	15.49	264	195796	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.46	112	297236	46.46	ng	0.00
7) Phenol-d6	6.47	99	375267	47.55	ng	-0.01
23) Nitrobenzene-d5	7.40	82	218619	34.49	ng	-0.02
41) 2,4,6-Tribromophenol	10.67	330	49114	34.43	ng	-0.01
44) 2-Fluorobiphenyl	9.20	172	428075	41.00	ng	-0.01
78) Terphenyl-d14	12.95	244	271937	32.82	ng	0.00
Target Compounds						
10) Phenol	6.49	94	22156	2.510	ng	95
49) Dimethylphthalate	9.59	163	61008	4.637	ng	100
74) Fluoranthene	12.59	202	35527	2.408	ng	96
77) Pyrene	12.82	202	31592	2.199	ng	98
87) Benzo(b)fluoranthene	15.06	252	26926m	2.237	ng	

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

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