

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101517\
 Data File : BF099515.D
 Acq On : 15 Oct 2017 15:54
 Operator : SJ/JU
 Sample : I5752-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 ENV-119-3(0-5)

Manual Integrations
 APPROVED

Sohil
 10/16/2017 7:45:31 PM

Quant Time: Oct 16 02:03:46 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	6.85	152	105086	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	419578	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	170930	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	241520	20.00	ng	0.00
75) Chrysene-d12	14.01	240	183115	20.00	ng	0.00
86) Perylene-d12	15.49	264	195020	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	707071	107.11	ng	0.00
7) Phenol-d6	6.49	99	875257	107.48	ng	0.00
23) Nitrobenzene-d5	7.41	82	522124	79.80	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	114543	81.89	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	888278	86.76	ng	-0.01
78) Terphenyl-d14	12.95	244	513149	63.73	ng	0.00
Target Compounds						
10) Phenol	6.49	94	55915	6.139	ng	96
49) Dimethylphthalate	9.59	163	164091	12.719	ng	99
70) Phenanthrene	11.39	178	30649	2.371	ng	97
74) Fluoranthene	12.58	202	53787	4.109	ng	100
77) Pyrene	12.82	202	53700	3.846	ng	97
80) Benzo(a)anthracene	14.00	228	29180	2.632	ng	99
82) Chrysene	14.04	228	26884	2.547	ng	98
85) Indeno(1,2,3-cd)pyrene	16.96	276	17073	2.022	ng	99
87) Benzo(b)fluoranthene	15.06	252	41095m	3.427	ng	
89) Benzo(a)pyrene	15.42	252	30675	2.787	ng	95

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

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