

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101517\
 Data File : BF099520.D
 Acq On : 15 Oct 2017 18:17
 Operator : SJ/JU
 Sample : I5752-10
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 TEST-PIT-116-C(0-5)

Manual Integrations
 APPROVED

Sohil
 10/16/2017 7:45:36 PM

Quant Time: Oct 16 02:26:54 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	95831	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	381449	20.00	ng	0.00
38) Acenaphthene-d10	9.88	164	153388	20.00	ng	-0.01
63) Phenanthrene-d10	11.37	188	220156	20.00	ng	0.00
75) Chrysene-d12	14.01	240	174816	20.00	ng	0.00
86) Perylene-d12	15.49	264	168315	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.47	112	686218	113.99	ng	0.00
7) Phenol-d6	6.48	99	813063	109.49	ng	0.00
23) Nitrobenzene-d5	7.41	82	473799	79.66	ng	-0.01
41) 2,4,6-Tribromophenol	10.67	330	128448	102.34	ng	0.00
44) 2-Fluorobiphenyl	9.20	172	799530	87.02	ng	-0.01
78) Terphenyl-d14	12.95	244	499992	65.04	ng	0.00
Target Compounds						
10) Phenol	6.49	94	48543	5.845	ng	97
49) Dimethylphthalate	9.59	163	114343	9.877	ng	99
70) Phenanthrene	11.39	178	46762	3.968	ng	98
74) Fluoranthene	12.59	202	69811	5.850	ng	97
77) Pyrene	12.82	202	65221	4.893	ng	97
80) Benzo(a)anthracene	14.00	228	33565	3.171	ng	97
82) Chrysene	14.04	228	31332	3.109	ng	98
87) Benzo(b)fluoranthene	15.06	252	37545m	3.628	ng	
89) Benzo(a)pyrene	15.42	252	28650	3.016	ng	# 94

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

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