

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099418.D
 Acq On : 13 Oct 2017 2:06
 Operator : SJ/JU
 Sample : I5779-03
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 FK-SF-03-003-B

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:06:41 PM

Quant Time: Oct 13 05:26:42 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	119844	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	456906	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	171575	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	255790	20.00	ng	0.00
75) Chrysene-d12	14.02	240	211699	20.00	ng	0.00
86) Perylene-d12	15.50	264	144702	20.00	ng	0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	822644	109.27	ng	0.01
7) Phenol-d6	6.49	99	960952	103.47	ng	0.00
23) Nitrobenzene-d5	7.42	82	572331	80.33	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	139460	99.33	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	922111	89.72	ng	0.00
78) Terphenyl-d14	12.96	244	651910	70.03	ng	0.00
Target Compounds						
10) Phenol	6.50	94	56913	5.480	ng	92
49) Dimethylphthalate	9.60	163	151364	11.689	ng	100
70) Phenanthrene	11.40	178	53444	3.903	ng	98
74) Fluoranthene	12.59	202	82748	5.968	ng	96
77) Pyrene	12.82	202	82480	5.109	ng	99
80) Benzo(a)anthracene	14.01	228	33084	2.581	ng	# 92
82) Chrysene	14.04	228	33734	2.764	ng	# 93
87) Benzo(b)fluoranthene	15.06	252	33206m	3.732	ng	
89) Benzo(a)pyrene	15.43	252	20776	2.544	ng	# 70

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

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