

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF101217\
 Data File : BF099417.D
 Acq On : 13 Oct 2017 1:38
 Operator : SJ/JU
 Sample : I5779-01
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

Sohil
 10/13/2017 3:06:40 PM

Quant Time: Oct 13 05:06:20 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	116864	20.00	ng	0.00
21) Naphthalene-d8	8.13	136	444632	20.00	ng	0.00
38) Acenaphthene-d10	9.89	164	162079	20.00	ng	0.00
63) Phenanthrene-d10	11.37	188	232143	20.00	ng	0.00
75) Chrysene-d12	14.02	240	216534	20.00	ng	0.00
86) Perylene-d12	15.49	264	165670	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.48	112	779132	106.13	ng	0.01
7) Phenol-d6	6.49	99	909004	100.37	ng	0.00
23) Nitrobenzene-d5	7.42	82	534516	77.09	ng	0.00
41) 2,4,6-Tribromophenol	10.68	330	127067	95.81	ng	0.00
44) 2-Fluorobiphenyl	9.21	172	896643	92.35	ng	0.00
78) Terphenyl-d14	12.96	244	593632	62.35	ng	0.00
Target Compounds						
10) Phenol	6.50	94	56496	5.578	ng	90
49) Dimethylphthalate	9.60	163	163187	13.340	ng	99
70) Phenanthrene	11.40	178	56785	4.570	ng	99
74) Fluoranthene	12.59	202	89882	7.143	ng	97
77) Pyrene	12.82	202	81780	4.953	ng	98
80) Benzo(a)anthracene	14.00	228	44547	3.398	ng	98
82) Chrysene	14.04	228	38812	3.109	ng	98
87) Benzo(b)fluoranthene	15.06	252	38995m	3.828	ng	
89) Benzo(a)pyrene	15.43	252	25800	2.759	ng	# 84
91) Benzo(g,h,i)perylene	17.42	276	14977	2.025	ng	# 87

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

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