

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100104.D
 Acq On : 3 Nov 2017 00:00
 Operator : SJ/JU
 Sample : I6247-01
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 S13-SP3-C

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:20:37 AM

Quant Time: Nov 03 04:10:36 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102317.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 01 18:13:08 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.77	152	82503	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	333221	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	143372	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	216173	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	137811	20.00	ng	-0.02
86) Perylene-d12	15.42	264	136111	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	451810	85.98	ng	0.00
7) Phenol-d6	6.42	99	523320	83.99	ng	-0.01
23) Nitrobenzene-d5	7.34	82	297047	57.39	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	100160	76.66	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	437304	47.06	ng	-0.01
78) Terphenyl-d14	12.88	244	261466	41.46	ng	-0.02
Target Compounds						
10) Phenol	6.43	94	25520	3.730	ng	86
49) Dimethylphthalate	9.52	163	130233	11.470	ng	100
70) Phenanthrene	11.32	178	103954	8.521	ng	97
74) Fluoranthene	12.52	202	133366	10.519	ng	96
77) Pyrene	12.75	202	120077	11.459	ng	97
80) Benzo(a)anthracene	13.94	228	69786	7.988	ng	98
82) Chrysene	13.98	228	69993	8.522	ng	97
85) Indeno(1,2,3-cd)pyrene	16.89	276	44888	5.953	ng	95
87) Benzo(b)fluoranthene	14.99	252	109475m	12.737	ng	
88) Benzo(k)fluoranthene	15.02	252	35191m	4.342	ng	
89) Benzo(a)pyrene	15.36	252	70898	9.183	ng	# 91
91) Benzo(a,h,i)perylene	17.34	276	43951	6.548	ng	96

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

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