

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100512.D
 Acq On : 13 Nov 2017 20:01
 Operator : SJ/JU
 Sample : I6301-15 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-02-110917-H

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:34 PM

Quant Time: Nov 14 03:06:00 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 13 14:08:38 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	77500	20.00	ng	0.00
21) Naphthalene-d8	8.17	136	284808	20.00	ng	0.00
38) Acenaphthene-d10	9.93	164	117977	20.00	ng	0.00
63) Phenanthrene-d10	11.42	188	181398	20.00	ng	0.00
75) Chrysene-d12	14.06	240	160832	20.00	ng	0.00
86) Perylene-d12	15.53	264	119652	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.50	112	111397	23.04	ng	0.00
7) Phenol-d6	6.50	99	133347	22.37	ng	-0.01
23) Nitrobenzene-d5	7.45	82	75714	17.58	ng	-0.01
41) 2,4,6-Tribromophenol	10.72	330	26810	22.30	ng	0.00
44) 2-Fluorobiphenyl	9.25	172	152113	19.63	ng	0.00
78) Terphenyl-d14	13.00	244	103637	14.81	ng	0.00
Target Compounds						
49) Dimethylphthalate	9.64	163	26636	2.957	ng	98
70) Phenanthrene	11.44	178	73061	7.650	ng	98
71) Anthracene	11.49	178	22663	2.339	ng	98
74) Fluoranthene	12.63	202	126174	12.465	ng	93
77) Pyrene	12.86	202	115595	10.083	ng	99
80) Benzo(a)anthracene	14.04	228	57250	5.911	ng	96
82) Chrysene	14.08	228	51140	5.453	ng	96
85) Indeno(1,2,3-cd)pyrene	17.02	276	18538	2.444	ng	# 86
87) Benzo(b)fluoranthene	15.10	252	46422m	6.549	ng	
88) Benzo(k)fluoranthene	15.13	252	17864m	2.667	ng	
89) Benzo(a)pyrene	15.47	252	31791	5.059	ng	# 89
91) Benzo(a,h,i)perylene	17.47	276	18742	3.725	ng	# 89

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

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