

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF111317\
 Data File : BF100523.D
 Acq On : 14 Nov 2017 4:33
 Operator : SJ/JU
 Sample : I6301-01
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

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

Manual Integrations
 APPROVED

SOHIL
 11/14/2017 5:20:44 PM

Quant Time: Nov 14 08:49:05 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 14 02:37:14 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.89	152	133074	20.00	ng	-0.01
21) Naphthalene-d8	8.17	136	501033	20.00	ng	-0.01
38) Acenaphthene-d10	9.93	164	198720	20.00	ng	-0.01
63) Phenanthrene-d10	11.41	188	300181	20.00	ng	-0.01
75) Chrysene-d12	14.05	240	277867	20.00	ng	-0.02
86) Perylene-d12	15.53	264	209664	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.51	112	777026	93.60	ng	0.00
7) Phenol-d6	6.52	99	937959	91.62	ng	0.00
23) Nitrobenzene-d5	7.45	82	583950	77.05	ng	-0.02
41) 2,4,6-Tribromophenol	10.72	330	168743	83.33	ng	-0.01
44) 2-Fluorobiphenyl	9.25	172	910915	69.80	ng	-0.02
78) Terphenyl-d14	13.00	244	608495	50.32	ng	-0.01
Target Compounds						
10) Phenol	6.53	94	40778	3.794	ng	97
49) Dimethylphthalate	9.63	163	175238	11.550	ng	99
74) Fluoranthene	12.62	202	77047	4.600	ng	97
77) Pyrene	12.85	202	72459	3.658	ng	98
80) Benzo(a)anthracene	14.04	228	48927m	2.924	ng	
82) Chrysene	14.07	228	43769	2.701	ng	97
87) Benzo(b)fluoranthene	15.09	252	49398m	3.977	ng	
89) Benzo(a)pyrene	15.46	252	33117	3.007	ng	# 91
91) Benzo(g,h,i)perylene	17.46	276	18585	2.108	ng	# 89

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

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