

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF110217\
 Data File : BF100100.D
 Acq On : 2 Nov 2017 22:11
 Operator : SJ/JU
 Sample : I6079-14 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 Q3-COMP

Manual Integrations
 APPROVED

Sohil
 11/3/2017 11:20:17 AM

Quant Time: Nov 03 03:07:20 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	65074	20.00	ng	0.00
21) Naphthalene-d8	8.05	136	258520	20.00	ng	-0.01
38) Acenaphthene-d10	9.80	164	113629	20.00	ng	-0.01
63) Phenanthrene-d10	11.30	188	179249	20.00	ng	-0.01
75) Chrysene-d12	13.94	240	113773	20.00	ng	-0.03
86) Perylene-d12	15.41	264	116991	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.40	112	111630	26.93	ng	0.00
7) Phenol-d6	6.42	99	147486	30.01	ng	0.00
23) Nitrobenzene-d5	7.34	82	79014	19.68	ng	-0.01
41) 2,4,6-Tribromophenol	10.60	330	27236	26.30	ng	-0.01
44) 2-Fluorobiphenyl	9.12	172	173082	23.50	ng	-0.01
78) Terphenyl-d14	12.87	244	102442	19.68	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.52	163	44626	4.959	ng	98
70) Phenanthrene	11.32	178	100654	9.950	ng	98
71) Anthracene	11.38	178	24948	2.430	ng	95
74) Fluoranthene	12.52	202	112052	10.659	ng	97
77) Pyrene	12.74	202	95041	10.986	ng	98
80) Benzo(a)anthracene	13.94	228	45529	6.313	ng	97
82) Chrysene	13.97	228	36471	5.379	ng	97
85) Indeno(1,2,3-cd)pyrene	16.89	276	18725	3.008	ng	97
87) Benzo(b)fluoranthene	14.99	252	44751m	6.058	ng	
88) Benzo(k)fluoranthene	15.02	252	18892m	2.712	ng	
89) Benzo(a)pyrene	15.35	252	33763	5.088	ng	# 95
91) Benzo(a,h,i)perylene	17.33	276	15117	2.620	ng	# 91

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

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