

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090717\
 Data File : BF098331.D
 Acq On : 7 Sep 2017 23:26
 Operator : SJ/JU
 Sample : I5071-04 5X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G61A-1.5-3

Manual Integrations
 APPROVED

Sohil
 9/8/2017 5:45:30 PM

Quant Time: Sep 08 03:04:16 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 05 17:28:20 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.69	152	107370	20.00	ng	-0.02
21) Naphthalene-d8	7.97	136	397428	20.00	ng	-0.01
38) Acenaphthene-d10	9.73	164	165399	20.00	ng	-0.01
63) Phenanthrene-d10	11.21	188	292914	20.00	ng	-0.01
75) Chrysene-d12	13.85	240	257868	20.00	ng	-0.02
86) Perylene-d12	15.26	264	173288	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	136004	19.27	ng	-0.02
7) Phenol-d6	6.33	99	186166	19.41	ng	-0.02
23) Nitrobenzene-d5	7.25	82	105782	14.46	ng	-0.02
41) 2,4,6-Tribromophenol	10.52	330	22969	16.01	ng	-0.02
44) 2-Fluorobiphenyl	9.05	172	172676	15.34	ng	-0.02
78) Terphenyl-d14	12.79	244	128003	10.35	ng	-0.02
Target Compounds						
49) Dimethylphthalate	9.45	163	26561	2.197	ng	99
74) Fluoranthene	12.42	202	73311	4.500	ng	96
77) Pyrene	12.65	202	69247	3.283	ng	97
80) Benzo(a)anthracene	13.83	228	32851	2.126	ng	98
82) Chrysene	13.87	228	36054m	2.345	ng	
87) Benzo(b)fluoranthene	14.85	252	33706m	3.086	ng	
89) Benzo(a)pyrene	15.19	252	20553	2.125	ng	# 86

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

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