

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090817\
 Data File : BF098359.D
 Acq On : 8 Sep 2017 19:15
 Operator : SJ/JU
 Sample : I5071-07
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 G58A-0-2.5

Manual Integrations
 APPROVED

Sohil
 9/11/2017 4:47:41 PM

Quant Time: Sep 09 01:44:15 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	6.68	152	100426	20.00	ng	-0.02
21) Naphthalene-d8	7.96	136	361884	20.00	ng	-0.02
38) Acenaphthene-d10	9.72	164	134661	20.00	ng	-0.02
63) Phenanthrene-d10	11.20	188	217687	20.00	ng	-0.02
75) Chrysene-d12	13.83	240	215227	20.00	ng	-0.03
86) Perylene-d12	15.24	264	163309	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.29	112	318334	48.22	ng	-0.02
7) Phenol-d6	6.32	99	465090	51.85	ng	-0.03
23) Nitrobenzene-d5	7.25	82	297623	44.68	ng	-0.03
41) 2,4,6-Tribromophenol	10.50	330	37342	31.96	ng	-0.03
44) 2-Fluorobiphenyl	9.04	172	447379	48.81	ng	-0.03
78) Terphenyl-d14	12.78	244	310617	30.10	ng	-0.03
Target Compounds						
49) Dimethylphthalate	9.44	163	63791	6.481	ng	Qvalue 100
70) Phenanthrene	11.22	178	44463	3.833	ng	99
74) Fluoranthene	12.41	202	68461	5.654	ng	98
77) Pyrene	12.63	202	76006	4.318	ng	98
80) Benzo(a)anthracene	13.82	228	41742	3.236	ng	96
82) Chrysene	13.86	228	37497m	2.922	ng	
87) Benzo(b)fluoranthene	14.84	252	41045m	3.987	ng	
89) Benzo(a)pyrene	15.19	252	26095	2.864	ng	# 75

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

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