

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090617\
 Data File : BF098287.D
 Acq On : 6 Sep 2017 22:02
 Operator : SJ/JU
 Sample : I5093-07
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 AU-06-090117-C

Manual Integrations
 APPROVED

Sohil
 9/7/2017 5:44:37 PM

Quant Time: Sep 07 06:28:50 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	6.70	152	99211	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	382286	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	152513	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	242273	20.00	ng	0.00
75) Chrysene-d12	13.84	240	223533	20.00	ng	-0.02
86) Perylene-d12	15.26	264	163209	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	649999	99.66	ng	0.00
7) Phenol-d6	6.35	99	872361	98.44	ng	0.00
23) Nitrobenzene-d5	7.26	82	523714	74.42	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	124316	93.94	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	757939	73.01	ng	-0.01
78) Terphenyl-d14	12.80	244	520439	48.55	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	32612	2.925	ng	99
70) Phenanthrene	11.23	178	40118	3.108	ng	96
74) Fluoranthene	12.42	202	93146	6.912	ng	99
77) Pyrene	12.65	202	103098	5.639	ng	98
80) Benzo(a)anthracene	13.84	228	54266	4.051	ng	95
82) Chrysene	13.87	228	50522	3.791	ng	97
87) Benzo(b)fluoranthene	14.86	252	52237m	5.078	ng	
89) Benzo(a)pyrene	15.20	252	34478	3.786	ng	94
91) Benzo(g,h,i)perylene	17.02	276	19554	2.528	ng	97

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

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