

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA F\DATA\BF090617\
 Data File : BF098293.D
 Acq On : 7 Sep 2017 00:51
 Operator : SJ/JU
 Sample : I5075-10 2X
 Misc :
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 292

Manual Integrations
 APPROVED

Sohil
 9/7/2017 5:44:41 PM

Quant Time: Sep 07 07:04:21 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	116684	20.00	ng	0.00
21) Naphthalene-d8	7.98	136	435367	20.00	ng	0.00
38) Acenaphthene-d10	9.73	164	165765	20.00	ng	0.00
63) Phenanthrene-d10	11.22	188	269974	20.00	ng	0.00
75) Chrysene-d12	13.85	240	242332	20.00	ng	-0.01
86) Perylene-d12	15.26	264	168122	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.30	112	396351	51.67	ng	0.00
7) Phenol-d6	6.34	99	524040	50.28	ng	-0.01
23) Nitrobenzene-d5	7.26	82	299518	37.37	ng	-0.01
41) 2,4,6-Tribromophenol	10.52	330	64447	44.81	ng	-0.01
44) 2-Fluorobiphenyl	9.06	172	468626	41.54	ng	-0.01
78) Terphenyl-d14	12.80	244	318316	27.39	ng	-0.01
Target Compounds						
49) Dimethylphthalate	9.45	163	69461	5.733	ng	99
70) Phenanthrene	11.24	178	69007	4.797	ng	98
74) Fluoranthene	12.42	202	146755	9.773	ng	98
77) Pyrene	12.65	202	139567	7.042	ng	97
80) Benzo(a)anthracene	13.84	228	72545	4.995	ng	97
82) Chrysene	13.87	228	64770	4.483	ng	97
85) Indeno(1,2,3-cd)pyrene	16.62	276	24419	2.298	ng	95
87) Benzo(b)fluoranthene	14.86	252	68058m	6.422	ng	
88) Benzo(k)fluoranthene	14.88	252	25060m	2.517	ng	
89) Benzo(a)pyrene	15.20	252	44800	4.775	ng	# 92
91) Benzo(g,h,i)perylene	17.03	276	26630	3.342	ng	95

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

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