

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA E\DATA\BE120517\
 Data File : BE094930.D
 Acq On : 6 Dec 2017 00:23
 Operator : SJ/JU
 Sample : I6698-11
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 FT-MW02S-20171130

Manual Integrations
 APPROVED

Sohil
 12/6/2017 5:10:30 PM

Quant Time: Dec 06 15:55:50 2017

Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE113017.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Fri Dec 01 13:33:28 2017

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.48	152	9999	0.40	ng	0.00
7) Naphthalene-d8	11.33	136	24276	0.40	ng	0.00
13) Acenaphthene-d10	15.08	164	33839	0.40	ng	0.00
19) Phenanthrene-d10	17.77	188	26739	0.40	ng	0.00
27) Chrysene-d12	21.89	240	24039m	0.40	ng	0.00
34) Perylene-d12	24.57	264	21936	0.40	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.95	112	3288	0.21	ng	0.00
5) Phenol-d6	7.60	99	2509	0.11	ng	0.00
8) Nitrobenzene-d5	9.66	82	7899	0.55	ng	0.00
11) 2-Methylnaphthalene-d10	12.86	152	16538	0.43	ng	0.00
14) 2,4,6-Tribromophenol	16.55	330	1725	0.30	ng	0.02
15) 2-Fluorobiphenyl	13.71	172	22779	0.16	ng	0.00
25) Fluoranthene-d10	19.76	212	128645	0.35	ng	0.00
29) Terphenyl-d14	20.33	244	22114	0.50	ng	0.00
Target Compounds						
28) Pyrene	20.14	202	10664	0.128	ng	Qvalue 97
30) Benzo(a)anthracene	21.87	228	3178	0.043	ng	# 87
31) Chrysene	21.92	228	3705m	0.045	ng	
32) Bis(2-ethylhexyl)phthalate	21.76	149	60987	0.499	ng	# 92
33) Indeno(1,2,3-cd)pyrene	27.43	276	3981	0.060	ng	# 85
35) Benzo(b)fluoranthene	23.73	252	6024m	0.075	ng	
36) Benzo(k)fluoranthene	23.79	252	3392	0.042	ng	87
39) Benzo(g,h,i)perylene	28.33	276	3792	0.059	ng	# 69

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

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