

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095438.D
 Acq On : 3 Feb 2018 15:36
 Operator : SJ/JU
 Sample : J1314-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C46

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:57 PM

Quant Time: Feb 05 03:58:56 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 03:32:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1779	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5470	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3630	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7588	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7177	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6438m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3113	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7696	0.27	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	16397	11.157	ng/ul	98
12) Phenanthrene	17.75	178	3200	0.125	ng/ul#	92
13) Anthracene	17.85	178	2336	0.100	ng/ul#	91
15) Fluoranthene	19.72	202	15176	0.459	ng/ul	83
17) Pyrene	20.08	202	14433	0.599	ng/ul#	78
18) Benzo(a)anthracene	21.78	228	4825	0.215	ng/ul#	88
19) Chrysene	21.84	228	5065	0.217	ng/ul#	88
21) Benzo(b)fluoranthene	23.62	252	9419m	0.415	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	2464m	0.119	ng/ul	
23) Benzo(a)pyrene	24.32	252	2445m	0.121	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	2855m	0.128	ng/ul	
25) Dibenzo(a,h)anthracene	27.25	278	599m	0.033	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	2505m	0.133	ng/ul	

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

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