

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095502.D
 Acq On : 5 Feb 2018 21:57
 Operator : SJ/JU
 Sample : J1314-06
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C43

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:31:08 PM

Quant Time: Feb 06 03:15:24 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 02:09:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1852	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5344	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3225	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7689	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7559	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7890	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	740	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	2749	0.09	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	1252	0.081	ng/ul#	66
11) Pentachlorophenol	17.36	266	19871	13.343	ng/ul	98
12) Phenanthrene	17.75	178	5215m	0.202	ng/ul	
13) Anthracene	17.85	178	6005m	0.254	ng/ul	
15) Fluoranthene	19.72	202	29843	0.891	ng/ul	82
17) Pyrene	20.07	202	34484	1.360	ng/ul#	79
18) Benzo(a)anthracene	21.78	228	10512	0.445	ng/ul#	89
19) Chrysene	21.84	228	11212	0.457	ng/ul#	90
21) Benzo(b)fluoranthene	23.62	252	24668m	0.886	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	6035m	0.237	ng/ul	
23) Benzo(a)pyrene	24.31	252	6682	0.269	ng/ul	92
24) Indeno(1,2,3-cd)pyrene	27.24	276	8720	0.320	ng/ul#	75
25) Dibenzo(a,h)anthracene	27.25	278	1992	0.089	ng/ul#	69
26) Benzo(g,h,i)perylene	28.11	276	9271	0.402	ng/ul	95

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

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