

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095250.D
 Acq On : 29 Jan 2018 17:08
 Operator : SJ/JU
 Sample : J1233-09
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B13

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:32:07 PM

Quant Time: Jan 30 05:53:26 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 Jan 30 04:43:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2371	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7287	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4809	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9799m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9754	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8884m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4375	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12862	0.34	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	2183	1.150	ng/ul	95
12) Phenanthrene	17.75	178	1199	0.036	ng/ul#	90
13) Anthracene	17.85	178	646	0.021	ng/ul#	66
15) Fluoranthene	19.72	202	959	0.022	ng/ul	86
17) Pyrene	20.08	202	877m	0.027	ng/ul	
21) Benzo(b)fluoranthene	23.63	252	1392m	0.044	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	599m	0.023	ng/ul	

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

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