

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095511.D
 Acq On : 6 Feb 2018 3:20
 Operator : SJ/JU
 Sample : J1315-17
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C80

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:31:38 PM

Quant Time: Feb 06 06:03:26 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 Feb 06 04:22:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1873	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5376	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3377	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7888	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7783	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7096m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2556	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7574	0.25	ng/ul	0.00
Target Compounds						
7) Acenaphthylene	14.74	152	917	0.056	ng/ul#	78
11) Pentachlorophenol	17.36	266	310	0.203	ng/ul	90
15) Fluoranthene	19.72	202	765	0.022	ng/ul	84
17) Pyrene	20.08	202	1319m	0.051	ng/ul	
21) Benzo(b)fluoranthene	23.62	252	885m	0.035	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1483m	0.060	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	1588m	0.077	ng/ul	

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

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