

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095517.D
 Acq On : 6 Feb 2018 6:52
 Operator : SJ/JU
 Sample : J1315-16
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C79

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:32:39 PM

Quant Time: Feb 06 08:04:39 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	8.44	152	1868	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5411	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3432	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7963	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7548	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7240m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3220	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	9348	0.31	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.75	152	1539	0.093	ng/ul	94
11) Pentachlorophenol	17.36	266	3250	2.107	ng/ul	98
12) Phenanthrene	17.75	178	687m	0.026	ng/ul	
13) Anthracene	17.85	178	799m	0.033	ng/ul	
15) Fluoranthene	19.72	202	1765	0.051	ng/ul	85
17) Pyrene	20.08	202	2643m	0.104	ng/ul	
19) Chrysene	21.84	228	922	0.038	ng/ul#	50
21) Benzo(b)fluoranthene	23.62	252	1305m	0.051	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.23	276	779m	0.031	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	1969m	0.093	ng/ul	

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

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