

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095470.D
 Acq On : 4 Feb 2018 10:39
 Operator : SJ/JU
 Sample : J1315-16
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C79

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:09 PM

Quant Time: Feb 05 07:28:46 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 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1851	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5602	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3802	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8245m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7612	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6929m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	4017	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	10740	0.34	ng/ul	0.00
Target Compounds					Qvalue	
7) Acenaphthylene	14.75	152	2375	0.130	ng/ul	94
11) Pentachlorophenol	17.36	266	4072	2.550	ng/ul	98
12) Phenanthrene	17.75	178	816	0.029	ng/ul#	80
13) Anthracene	17.85	178	1152m	0.045	ng/ul	
15) Fluoranthene	19.72	202	1935	0.054	ng/ul	84
17) Pyrene	20.08	202	2602m	0.102	ng/ul	
18) Benzo(a)anthracene	21.78	228	755	0.032	ng/ul#	45
19) Chrysene	21.84	228	994	0.040	ng/ul#	53
21) Benzo(b)fluoranthene	23.62	252	1495m	0.061	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	896m	0.037	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	2183m	0.108	ng/ul	

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

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