

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095469.D
 Acq On : 4 Feb 2018 10:03
 Operator : SJ/JU
 Sample : J1315-18
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C81

Manual Integrations
 APPROVED

Sohil
 2/5/2018 7:01:07 PM

Quant Time: Feb 05 07:21:10 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	1956	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	5895	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3977	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8623	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	8225	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	7711m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3297	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8875	0.27	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	1158	0.061	ng/ul#	78
11) Pentachlorophenol	17.36	266	773	0.463	ng/ul	98
12) Phenanthrene	17.75	178	832	0.029	ng/ul#	64
13) Anthracene	17.85	178	1428m	0.054	ng/ul	
15) Fluoranthene	19.72	202	2972	0.079	ng/ul	82
17) Pyrene	20.08	202	5138m	0.186	ng/ul	
18) Benzo(a)anthracene	21.79	228	2893	0.113	ng/ul#	65
19) Chrysene	21.84	228	3203	0.120	ng/ul#	78
21) Benzo(b)fluoranthene	23.62	252	4514m	0.166	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	1212m	0.049	ng/ul	
23) Benzo(a)pyrene	24.32	252	1987m	0.082	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1438m	0.054	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	1152m	0.051	ng/ul	

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

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