

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095518.D
 Acq On : 6 Feb 2018 7:27
 Operator : SJ/JU
 Sample : J1315-18
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6C81

Manual Integrations
 APPROVED

Sohil
 2/6/2018 2:32:41 PM

Quant Time: Feb 06 08:12:44 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	1723	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	4861	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3207	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7442	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7034	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6608m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2450	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7350	0.26	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.75	152	667	0.043	ng/ul#	74
11) Pentachlorophenol	17.36	266	563	0.391	ng/ul	93
12) Phenanthrene	17.75	178	614	0.025	ng/ul#	63
13) Anthracene	17.85	178	1095m	0.048	ng/ul	
15) Fluoranthene	19.72	202	2535	0.078	ng/ul	83
17) Pyrene	20.07	202	3897m	0.165	ng/ul	
18) Benzo(a)anthracene	21.78	228	3302	0.150	ng/ul#	47
19) Chrysene	21.84	228	2654	0.116	ng/ul#	76
21) Benzo(b)fluoranthene	23.62	252	3804m	0.163	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	993m	0.047	ng/ul	
23) Benzo(a)pyrene	24.31	252	1585m	0.076	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1138m	0.050	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	994m	0.051	ng/ul	

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

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