

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095270.D
 Acq On : 30 Jan 2018 5:04
 Operator : SJ/JU
 Sample : J1221-02
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6A23

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:33:03 PM

Quant Time: Jan 30 07:13:41 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 Jan 30 06:14:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2121	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6456	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4080	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8774m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9228	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8425m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4228	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	11775	0.35	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	14.74	152	639	0.033	ng/ul#	73
11) Pentachlorophenol	17.36	266	1360	0.800	ng/ul	93
12) Phenanthrene	17.75	178	983	0.033	ng/ul#	76
13) Anthracene	17.85	178	3865	0.143	ng/ul	94
15) Fluoranthene	19.73	202	2401	0.063	ng/ul	84
17) Pyrene	20.08	202	2538m	0.082	ng/ul	
18) Benzo(a)anthracene	21.79	228	1726	0.060	ng/ul#	76
19) Chrysene	21.82	228	1803	0.060	ng/ul#	74
21) Benzo(b)fluoranthene	23.63	252	2796m	0.094	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	668m	0.025	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1407m	0.048	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	897m	0.036	ng/ul	

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

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