

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095251.D
 Acq On : 29 Jan 2018 17:44
 Operator : SJ/JU
 Sample : J1233-19
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B23

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:32:08 PM

Quant Time: Jan 30 05:58:47 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 04:43:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2234	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6905	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4511	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	11912m	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	9775	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9359m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4325	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12621	0.28	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.31	128	102220	5.287	ng/ul#	87
5) 2-Methylnaphthalene	12.89	142	30384	2.353	ng/ul	100
8) Acenaphthene	15.08	153	3632	0.216	ng/ul	95
9) Fluorene	16.03	166	1862m	0.107	ng/ul	
11) Pentachlorophenol	17.36	266	249442	108.114	ng/ul	96
12) Phenanthrene	17.75	178	7127	0.178	ng/ul#	86
13) Anthracene	17.85	178	1431	0.039	ng/ul#	92
15) Fluoranthene	19.72	202	5554	0.107	ng/ul	84
17) Pyrene	20.08	202	3430m	0.105	ng/ul	
18) Benzo(a)anthracene	21.79	228	1357	0.044	ng/ul#	84
19) Chrysene	21.84	228	884	0.028	ng/ul#	80
21) Benzo(b)fluoranthene	23.62	252	868m	0.026	ng/ul	

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

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