

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095266.D
 Acq On : 30 Jan 2018 2:44
 Operator : SJ/JU
 Sample : J1244-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B47

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:32:56 PM

Quant Time: Jan 30 06:58:40 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	2548	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7550	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4769	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9797	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10271	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9155m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3743	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	9827	0.26	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.31	128	2430	0.115	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	1436	0.102	ng/ul	99
7) Acenaphthylene	14.74	152	543	0.024	ng/ul#	69
8) Acenaphthene	15.08	153	3325	0.187	ng/ul	93
9) Fluorene	16.03	166	2516	0.137	ng/ul#	79
11) Pentachlorophenol	17.36	266	1716	0.904	ng/ul	95
12) Phenanthrene	17.75	178	11017	0.334	ng/ul#	88
13) Anthracene	17.85	178	2447	0.081	ng/ul	95
15) Fluoranthene	19.73	202	4755	0.111	ng/ul	84
17) Pyrene	20.08	202	4074m	0.118	ng/ul	
18) Benzo(a)anthracene	21.79	228	1520	0.047	ng/ul#	87
19) Chrysene	21.84	228	1507	0.045	ng/ul#	79
21) Benzo(b)fluoranthene	23.63	252	2714m	0.084	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	878m	0.030	ng/ul	
23) Benzo(a)pyrene	24.32	252	941m	0.033	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	1351m	0.043	ng/ul	
26) Benzo(g,h,i)perylene	28.11	276	1503m	0.056	ng/ul	

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

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