

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095249.D
 Acq On : 29 Jan 2018 16:32
 Operator : SJ/JU
 Sample : J1233-10
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6B14

Manual Integrations
 APPROVED

Sohil
 1/30/2018 3:32:05 PM

Quant Time: Jan 30 05:22:12 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	2548	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7814	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5136	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10825m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10679	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9521m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4660	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12747	0.31	ng/ul	0.00
Target Compounds				Qvalue		
11) Pentachlorophenol	17.36	266	3329	1.588	ng/ul	97
12) Phenanthrene	17.75	178	1760	0.048	ng/ul#	92
13) Anthracene	17.85	178	911	0.027	ng/ul#	80
15) Fluoranthene	19.72	202	2794	0.059	ng/ul	84
17) Pyrene	20.08	202	3025m	0.084	ng/ul	
18) Benzo(a)anthracene	21.79	228	1019	0.031	ng/ul#	79
19) Chrysene	21.84	228	1170	0.034	ng/ul#	73
21) Benzo(b)fluoranthene	23.62	252	2396m	0.071	ng/ul	
22) Benzo(k)fluoranthene	23.65	252	975m	0.032	ng/ul	
23) Benzo(a)pyrene	24.32	252	820m	0.027	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	825m	0.025	ng/ul	
26) Benzo(g,h,i)perylene	28.12	276	1114m	0.040	ng/ul	

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

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