

Data Path : U:\HPCHEM1\BNA E\DATA\BE012018\
 Data File : BE095083.D
 Acq On : 20 Jan 2018 11:11
 Operator : SJ/JU
 Sample : J1216-05
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 B0AE8

Manual Integrations
 APPROVED

Sohil
 1/22/2018 5:22:24 PM

Quant Time: Jan 22 12:04:15 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE011918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 22 04:29:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4951	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	16147	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	10181	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	20830m	0.40	ng/ul	0.00
16) Chrysene-d12	21.83	240	22508	0.40	ng/ul	0.00
20) Perylene-d12	24.48	264	17514m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.84	152	9994	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	28563	0.44	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	14.77	152	3377	0.074	ng/ul	98
11) Pentachlorophenol	17.39	266	350	0.284	ng/ul	91
12) Phenanthrene	17.78	178	10635	0.192	ng/ul#	88
13) Anthracene	17.87	178	2204	0.046	ng/ul#	95
15) Fluoranthene	19.75	202	40121	0.568	ng/ul	84
17) Pyrene	20.10	202	39373	0.481	ng/ul#	83
18) Benzo(a)anthracene	21.81	228	24528	0.394	ng/ul#	86
19) Chrysene	21.87	228	19386	0.279	ng/ul	98
21) Benzo(b)fluoranthene	23.66	252	25003m	0.424	ng/ul	
22) Benzo(k)fluoranthene	23.71	252	9234m	0.170	ng/ul	
23) Benzo(a)pyrene	24.36	252	18073m	0.359	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.31	276	10411m	0.233	ng/ul	
25) Dibenzo(a,h)anthracene	27.33	278	3035m	0.095	ng/ul	
26) Benzo(g,h,i)perylene	28.19	276	7874m	0.177	ng/ul	

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

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