

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031418\
 Data File : BE095773.D
 Acq On : 14 Mar 2018 20:34
 Operator : SJ/JU
 Sample : J1813-16
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6F34

Manual Integrations
 APPROVED

Sohil
 3/15/2018 6:54:23 PM

Quant Time: Mar 15 12:10:16 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 03:09:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1885	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5495	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3383	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7368m	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7106m	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	7883	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	1638	0.18	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	4314m	0.18	ng/ul	0.00
Target Compounds					Ovalue	
8) Acenaphthene	15.07	153	512	0.040	ng/ul#	93
9) Fluorene	16.03	166	409	0.029	ng/ul#	83
12) Phenanthrene	17.75	178	8454m	0.383	ng/ul	
13) Anthracene	17.84	178	1675m	0.081	ng/ul	
15) Fluoranthene	19.72	202	21934m	0.821	ng/ul	
17) Pyrene	20.07	202	18600m	0.689	ng/ul	
18) Benzo(a)anthracene	21.78	228	12990	0.567	ng/ul#	87
19) Chrysene	21.84	228	10454	0.448	ng/ul	97
21) Benzo(b)fluoranthene	23.61	252	14712m	0.539	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	5239m	0.202	ng/ul	
23) Benzo(a)pyrene	24.30	252	10127m	0.406	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.22	276	6943	0.251	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.24	278	1685	0.076	ng/ul#	87
26) Benzo(g,h,i)perylene	28.10	276	6362	0.268	ng/ul	95

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

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