

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031418\
 Data File : BE095784.D
 Acq On : 15 Mar 2018 3:08
 Operator : SJ/JU
 Sample : J1813-17
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6F35

Manual Integrations
 APPROVED

Sohil
 3/15/2018 6:54:56 PM

Quant Time: Mar 15 05:42:21 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 05:35:28 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	2049	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	6320	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	4042	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8517m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7852m	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	7542m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3126	0.29	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7740m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	15.07	153	549	0.036	ng/ul	93
9) Fluorene	16.04	166	440	0.026	ng/ul#	90
12) Phenanthrene	17.75	178	10525m	0.413	ng/ul	
13) Anthracene	17.85	178	2023m	0.085	ng/ul	
15) Fluoranthene	19.72	202	22430m	0.726	ng/ul	
17) Pyrene	20.07	202	18171m	0.609	ng/ul	
18) Benzo(a)anthracene	21.78	228	8316m	0.328	ng/ul	
19) Chrysene	21.84	228	9581m	0.371	ng/ul	
21) Benzo(b)fluoranthene	23.61	252	12217m	0.468	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	4079m	0.164	ng/ul	
23) Benzo(a)pyrene	24.31	252	7879	0.330	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	27.23	276	5665	0.214	ng/ul#	72
25) Dibenzo(a,h)anthracene	27.24	278	1487	0.070	ng/ul#	44
26) Benzo(g,h,i)perylene	28.11	276	5449	0.240	ng/ul#	94

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

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