

Data Path : Z:\HPCHEM1\BNA_E\Data\BE031418\
 Data File : BE095785.D
 Acq On : 15 Mar 2018 3:44
 Operator : SJ/JU
 Sample : J1813-08
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 F6F29

Manual Integrations
 APPROVED

Sohil
 3/15/2018 6:54:59 PM

Quant Time: Mar 20 16:09:41 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	2033	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	6158	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3847	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7910m	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	7485m	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	8732	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3122	0.30	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	7518m	0.30	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.30	128	1512	0.088	ng/ul#	87
5) 2-Methylnaphthalene	12.88	142	1945	0.163	ng/ul	99
7) Acenaphthylene	14.74	152	2846	0.147	ng/ul	96
12) Phenanthrene	17.75	178	12583m	0.531	ng/ul	
13) Anthracene	17.85	178	1597m	0.072	ng/ul	
15) Fluoranthene	19.72	202	36180m	1.261	ng/ul	
17) Pyrene	20.07	202	40495m	1.425	ng/ul	
18) Benzo(a)anthracene	21.78	228	20513m	0.849	ng/ul	
19) Chrysene	21.84	228	25916m	1.054	ng/ul	
21) Benzo(b)fluoranthene	23.61	252	42272m	1.398	ng/ul	
22) Benzo(k)fluoranthene	23.66	252	11142m	0.388	ng/ul	
23) Benzo(a)pyrene	24.32	252	26273m	0.951	ng/ul	
24) Indeno(1,2,3-cd)pyrene	27.24	276	21247	0.694	ng/ul#	72
25) Dibenzo(a,h)anthracene	27.25	278	4796	0.196	ng/ul#	79
26) Benzo(g,h,i)perylene	28.12	276	20193	0.768	ng/ul	94

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

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