

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097169.D
 Acq On : 3 Aug 2018 2:22
 Operator : SJ/JU
 Sample : J4171-19
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JJ1L2

Manual Integrations
 APPROVED

Sohil
 8/3/2018 5:55:20 PM

Quant Time: Aug 03 07:35:53 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 02 23:40:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1322	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6907	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4461	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	11362m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	12359m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10672m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	2797	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9614m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.81	178	3881m	0.134	ng/ul	
13) Anthracene	16.90	178	831m	0.031	ng/ul	
15) Fluoranthene	18.84	202	6053m	0.158	ng/ul	
17) Pyrene	19.20	202	6643	0.201	ng/ul#	85
18) Benzo(a)anthracene	20.96	228	2506	0.071	ng/ul#	84
19) Chrysene	21.01	228	2319m	0.064	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	4953m	0.170	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1661m	0.053	ng/ul	
23) Benzo(a)pyrene	23.11	252	2696	0.093	ng/ul	94
24) Indeno(1,2,3-cd)pyrene	25.49	276	5731	0.157	ng/ul#	73
25) Dibenzo(a,h)anthracene	25.50	278	913	0.031	ng/ul#	52
26) Benzo(g,h,i)perylene	26.19	276	5781	0.184	ng/ul	96

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

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