

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097168.D
 Acq On : 3 Aug 2018 1:44
 Operator : SJ/JU
 Sample : J4171-14
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JJ1K5

Manual Integrations
 APPROVED

Sohil
 8/3/2018 5:55:19 PM

Quant Time: Aug 03 08:06:34 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	1189	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6202	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4046	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10349m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11141m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9482m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	3298	0.33	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	10907m	0.31	ng/ul	0.00
Target Compounds						
3) Naphthalene	10.18	128	1735	0.116	ng/ul#	94
5) 2-Methylnaphthalene	11.80	142	1474	0.149	ng/ul	98
12) Phenanthrene	16.81	178	2951m	0.112	ng/ul	
15) Fluoranthene	18.84	202	3581m	0.103	ng/ul	
17) Pyrene	19.20	202	4186m	0.140	ng/ul	
18) Benzo(a)anthracene	20.96	228	1815	0.057	ng/ul#	89
19) Chrysene	21.01	228	2192	0.067	ng/ul#	89
21) Benzo(b)fluoranthene	22.53	252	2599m	0.101	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	1088m	0.039	ng/ul	
23) Benzo(a)pyrene	23.11	252	1793	0.069	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.49	276	1432	0.044	ng/ul#	86
26) Benzo(g,h,i)perylene	26.19	276	1488	0.053	ng/ul#	79

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

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