

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097421.D
 Acq On : 7 Sep 2018 12:41
 Operator : SJ/JU
 Sample : J4688-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z24

Manual Integrations
 APPROVED

Sohil
 9/11/2018 2:44:49 PM

Quant Time: Sep 08 00:16:38 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 07 23:49:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	909	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4655	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3251	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8970m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10712m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9976m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1849	0.26	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	7680m	0.26	ng/ul	0.00
Target Compounds				Qvalue		
12) Phenanthrene	16.80	178	3314m	0.144	ng/ul	
13) Anthracene	16.89	178	411m	0.021	ng/ul	
15) Fluoranthene	18.83	202	6452m	0.218	ng/ul	
17) Pyrene	19.19	202	7304m	0.253	ng/ul	
18) Benzo(a)anthracene	20.95	228	2913m	0.106	ng/ul	
19) Chrysene	21.00	228	3365m	0.106	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	4098m	0.153	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	1731m	0.058	ng/ul	
23) Benzo(a)pyrene	23.10	252	2167	0.088	ng/ul#	55
24) Indeno(1,2,3-cd)pyrene	25.49	276	1750	0.055	ng/ul#	80
26) Benzo(g,h,i)perylene	26.19	276	1828	0.067	ng/ul#	87

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

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