

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097308.D
 Acq On : 17 Aug 2018 22:14
 Operator : SJ/JU
 Sample : J4424-16
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z63

Manual Integrations
 APPROVED

Sohil
 8/20/2018 4:15:37 PM

Quant Time: Aug 18 00:42:26 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 23:33:44 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1082	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	5546	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	3075	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7212m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	6650m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	5450m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1995	0.22	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	4845m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.80	178	1955m	0.105	ng/ul	
13) Anthracene	16.90	178	386m	0.025	ng/ul	
15) Fluoranthene	18.84	202	3362m	0.170	ng/ul	
17) Pyrene	19.20	202	3976m	0.198	ng/ul	
18) Benzo(a)anthracene	20.96	228	1483	0.090	ng/ul#	85
19) Chrysene	21.01	228	1992	0.101	ng/ul	94
21) Benzo(b)fluoranthene	22.54	252	2285m	0.155	ng/ul	
22) Benzo(k)fluoranthene	22.58	252	910m	0.050	ng/ul	
23) Benzo(a)pyrene	23.11	252	1519m	0.111	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.51	276	1094	0.065	ng/ul#	78
26) Benzo(g,h,i)perylene	26.21	276	1016	0.075	ng/ul#	82

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

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