

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE091218\
 Data File : BE097471.D
 Acq On : 13 Sep 2018 3:40
 Operator : SJ/JU
 Sample : J4792-19
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YZ5

Manual Integrations
 APPROVED

Sohil
 9/13/2018 5:42:38 PM

Quant Time: Sep 13 08:02:43 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 : Thu Sep 13 07:48:38 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.37	152	1306	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	7305	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	5237	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	11561m	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	10680m	0.40	ng/ul	0.02
20) Perylene-d12	23.24	264	10714m	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	3348	0.29	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	9707m	0.25	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.16	128	14764	0.829	ng/ul#	91
5) 2-Methylnaphthalene	11.79	142	3066	0.272	ng/ul	97
7) Acenaphthylene	13.72	152	3022	0.158	ng/ul#	71
8) Acenaphthene	14.07	153	2417	0.147	ng/ul#	88
9) Fluorene	15.07	166	4560	0.243	ng/ul#	82
12) Phenanthrene	16.80	178	44046m	1.480	ng/ul	
13) Anthracene	16.90	178	13304m	0.532	ng/ul	
15) Fluoranthene	18.84	202	106989m	2.798	ng/ul	
17) Pyrene	19.21	202	144156m	5.016	ng/ul	
18) Benzo(a)anthracene	20.97	228	82610m	3.011	ng/ul	
19) Chrysene	21.02	228	109988m	3.480	ng/ul	
21) Benzo(b)fluoranthene	22.56	252	134641m	4.693	ng/ul	
22) Benzo(k)fluoranthene	22.60	252	41264m	1.290	ng/ul	
23) Benzo(a)pyrene	23.14	252	70568m	2.659	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.55	276	56788m	1.651	ng/ul	
25) Dibenzo(a,h)anthracene	25.56	278	17186m	0.590	ng/ul	
26) Benzo(g,h,i)perylene	26.27	276	52930m	1.805	ng/ul	

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

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