

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097422.D
 Acq On : 7 Sep 2018 13:19
 Operator : SJ/JU
 Sample : J4688-15
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z21

Manual Integrations
 APPROVED

Sohil
 9/11/2018 2:44:03 PM

Quant Time: Sep 08 00:22:57 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	925	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4831	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	3230	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	8664m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	9981m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	9396m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	1360	0.18	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	5103m	0.18	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.16	128	2476	0.210	ng/ul#	96
5) 2-Methylnaphthalene	11.79	142	2247	0.302	ng/ul	98
7) Acenaphthylene	13.72	152	487	0.041	ng/ul#	57
9) Fluorene	15.07	166	276	0.024	ng/ul#	62
12) Phenanthrene	16.80	178	9877m	0.443	ng/ul	
13) Anthracene	16.89	178	1039m	0.055	ng/ul	
15) Fluoranthene	18.83	202	14914m	0.521	ng/ul	
17) Pyrene	19.20	202	16280	0.606	ng/ul#	79
18) Benzo(a)anthracene	20.95	228	7883m	0.307	ng/ul	
19) Chrysene	21.00	228	10071m	0.341	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	10657m	0.424	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	3599m	0.128	ng/ul	
23) Benzo(a)pyrene	23.10	252	7772m	0.334	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.48	276	5142m	0.170	ng/ul	
25) Dibenzo(a,h)anthracene	25.50	278	1400m	0.055	ng/ul	
26) Benzo(g,h,i)perylene	26.19	276	5409m	0.210	ng/ul	

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

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