

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE090718\
 Data File : BE097425.D
 Acq On : 7 Sep 2018 15:10
 Operator : SJ/JU
 Sample : J4688-06
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3YY1

Manual Integrations
 APPROVED

Sohil
 9/11/2018 2:42:37 PM

Quant Time: Sep 08 00:41:31 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	1054	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	5578	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	4108	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	10747m	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10856m	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	10697m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.71	152	1966	0.23	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	6609m	0.19	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.17	128	4712	0.346	ng/ul#	93
5) 2-Methylnaphthalene	11.79	142	1188	0.138	ng/ul	98
7) Acenaphthylene	13.72	152	1259	0.084	ng/ul#	55
8) Acenaphthene	14.07	153	1005	0.078	ng/ul#	90
9) Fluorene	15.07	166	1310	0.089	ng/ul#	81
12) Phenanthrene	16.80	178	16183m	0.585	ng/ul	
13) Anthracene	16.89	178	3831m	0.165	ng/ul	
15) Fluoranthene	18.83	202	29639m	0.834	ng/ul	
17) Pyrene	19.20	202	30590m	1.047	ng/ul	
18) Benzo(a)anthracene	20.95	228	15929m	0.571	ng/ul	
19) Chrysene	21.00	228	10875m	0.338	ng/ul	
21) Benzo(b)fluoranthene	22.53	252	57879m	2.020	ng/ul	
22) Benzo(k)fluoranthene	22.57	252	19422m	0.608	ng/ul	
23) Benzo(a)pyrene	23.10	252	42836m	1.616	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.48	276	41333m	1.203	ng/ul	
25) Dibenzo(a,h)anthracene	25.50	278	9724m	0.334	ng/ul	
26) Benzo(g,h,i)perylene	26.19	276	44868m	1.532	ng/ul	

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

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