

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097836.D
 Acq On : 10 Oct 2018 20:14
 Operator : SJ/JU
 Sample : J5184-04
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC23

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:33:37 PM

Quant Time: Oct 11 01:43:19 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE101018.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Oct 11 01:20:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	4830	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	23870	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	16569	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	39553m	0.40	ng/ul	0.00
16) Chrysene-d12	21.49	240	35995m	0.40	ng/ul	0.00
20) Perylene-d12	24.07	264	32420m	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	11490	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	32398m	0.25	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	1797	0.031	ng/ul#	83
5) 2-Methylnaphthalene	12.36	142	1340	0.035	ng/ul	99
7) Acenaphthylene	14.26	152	7028	0.120	ng/ul#	77
9) Fluorene	15.59	166	2661	0.045	ng/ul#	73
12) Phenanthrene	17.34	178	60320m	0.638	ng/ul	
13) Anthracene	17.43	178	15569m	0.194	ng/ul	
15) Fluoranthene	19.36	202	160736m	1.290	ng/ul	
17) Pyrene	19.72	202	157152m	1.777	ng/ul	
18) Benzo(a)anthracene	21.47	228	97660m	0.978	ng/ul	
19) Chrysene	21.53	228	342005	3.548	ng/ul#	86
21) Benzo(b)fluoranthene	23.28	252	212723m	2.539	ng/ul	
22) Benzo(k)fluoranthene	23.32	252	55710m	0.587	ng/ul	
23) Benzo(a)pyrene	23.96	252	108636	1.417	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	26.78	276	96155	1.105	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.80	278	25619m	0.352	ng/ul	
26) Benzo(a,h,i)perylene	27.62	276	122329	1.489	ng/ul#	63

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

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