

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097845.D
 Acq On : 11 Oct 2018 1:49
 Operator : SJ/JU
 Sample : J5183-08
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC47

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:33:59 PM

Quant Time: Oct 11 02:57:18 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	3484	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	18049	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	14059	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	37113m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	38201m	0.40	ng/ul	0.00
20) Perylene-d12	24.04	264	34863m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	9056	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	35145m	0.29	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.74	128	5823	0.134	ng/ul	96
5) 2-Methylnaphthalene	12.36	142	4539	0.157	ng/ul	100
7) Acenaphthylene	14.27	152	2620	0.053	ng/ul#	72
8) Acenaphthene	14.61	153	11400	0.276	ng/ul	98
9) Fluorene	15.60	166	14388	0.289	ng/ul	98
12) Phenanthrene	17.34	178	293201m	3.306	ng/ul	
13) Anthracene	17.44	178	58026m	0.771	ng/ul	
15) Fluoranthene	19.36	202	491094m	4.201	ng/ul	
17) Pyrene	19.72	202	475092m	5.062	ng/ul	
18) Benzo(a)anthracene	21.47	228	264485m	2.494	ng/ul	
19) Chrysene	21.52	228	565463	5.528	ng/ul	97
21) Benzo(b)fluoranthene	23.25	252	636260m	7.062	ng/ul	
22) Benzo(k)fluoranthene	23.30	252	165716m	1.623	ng/ul	
23) Benzo(a)pyrene	23.93	252	264312	3.206	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.73	276	317272	3.390	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	77454	0.990	ng/ul#	88
26) Benzo(a,h,i)perylene	27.55	276	367742	4.163	ng/ul	98

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

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