

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097840.D
 Acq On : 10 Oct 2018 22:43
 Operator : SJ/JU
 Sample : J5183-03
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC18

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:33:47 PM

Quant Time: Oct 11 02:11:26 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	2715	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	13386	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	10361	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	28474m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	34184m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	31644m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	5399	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	23909m	0.26	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.34	178	5623m	0.083	ng/ul	
15) Fluoranthene	19.36	202	27519m	0.307	ng/ul	
17) Pyrene	19.72	202	31282	0.372	ng/ul	96
18) Benzo(a)anthracene	21.47	228	29763m	0.314	ng/ul	
19) Chrysene	21.52	228	179357m	1.959	ng/ul	
21) Benzo(b)fluoranthene	23.25	252	153169m	1.873	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	36637m	0.395	ng/ul	
23) Benzo(a)pyrene	23.92	252	16487	0.220	ng/ul#	56
24) Indeno(1,2,3-cd)pyrene	26.72	276	44323	0.522	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.75	278	13045	0.184	ng/ul#	54
26) Benzo(g,h,i)perylene	27.54	276	49472	0.617	ng/ul#	93

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

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