

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097834.D
 Acq On : 10 Oct 2018 19:00
 Operator : SJ/JU
 Sample : J5184-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC35

Manual Integrations
 APPROVED

Sohil
 10/11/2018 5:33:32 PM

Quant Time: Oct 11 01:34:45 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	7.88	152	3147	0.40	ng/ul	0.00
2) Naphthalene-d8	10.68	136	15417	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	12600	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	35269m	0.40	ng/ul	0.00
16) Chrysene-d12	21.48	240	39201m	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	34907m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6874	0.28	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	31516m	0.27	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	2471	0.067	ng/ul#	87
8) Acenaphthene	14.61	153	1585	0.043	ng/ul	95
12) Phenanthrene	17.34	178	5440m	0.065	ng/ul	
15) Fluoranthene	19.36	202	9156m	0.082	ng/ul	
17) Pyrene	19.72	202	10621m	0.110	ng/ul	
18) Benzo(a)anthracene	21.46	228	3729m	0.034	ng/ul	
19) Chrysene	21.52	228	7652m	0.073	ng/ul	
21) Benzo(b)fluoranthene	23.25	252	9839m	0.109	ng/ul	
22) Benzo(k)fluoranthene	23.29	252	3466m	0.034	ng/ul	
23) Benzo(a)pyrene	23.91	252	3925	0.048	ng/ul#	31
24) Indeno(1,2,3-cd)pyrene	26.70	276	4299	0.046	ng/ul#	100
26) Benzo(g,h,i)perylene	27.54	276	5041	0.057	ng/ul#	84

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

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