

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097866.D
 Acq On : 11 Oct 2018 20:33
 Operator : SJ/JU
 Sample : J5183-02
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC17

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:34:56 PM

Quant Time: Oct 12 04:23:35 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE101118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 12 03:56:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2388	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	12212	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.55	164	8793	0.40	ng/ul	0.01
10) Phenanthrene-d10	17.31	188	20979m	0.40	ng/ul	0.01
16) Chrysene-d12	21.49	240	20081m	0.40	ng/ul	0.01
20) Perylene-d12	24.06	264	19220m	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	4627	0.22	ng/ul	0.00
14) Fluoranthene-d10	19.34	212	15483m	0.23	ng/ul	0.01
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	3875	0.130	ng/ul#	90
5) 2-Methylnaphthalene	12.37	142	2606	0.121	ng/ul	99
7) Acenaphthylene	14.27	152	1989	0.060	ng/ul#	79
8) Acenaphthene	14.61	153	1784	0.069	ng/ul	99
9) Fluorene	15.61	166	2472	0.073	ng/ul#	90
12) Phenanthrene	17.35	178	37361	0.752	ng/ul	95
13) Anthracene	17.44	178	6906m	0.150	ng/ul	
15) Fluoranthene	19.37	202	63552m	0.996	ng/ul	
17) Pyrene	19.73	202	72928m	1.219	ng/ul	
18) Benzo(a)anthracene	21.47	228	38978m	0.642	ng/ul	
19) Chrysene	21.53	228	55808m	1.023	ng/ul	
21) Benzo(b)fluoranthene	23.27	252	75386m	1.468	ng/ul	
22) Benzo(k)fluoranthene	23.32	252	22905m	0.417	ng/ul	
23) Benzo(a)pyrene	23.95	252	47523	0.942	ng/ul#	47
24) Indeno(1,2,3-cd)pyrene	26.77	276	38517	0.673	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.79	278	9796	0.207	ng/ul#	1
26) Benzo(g,h,i)perylene	27.60	276	42490	0.852	ng/ul#	72

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

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