

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097861.D
 Acq On : 11 Oct 2018 17:28
 Operator : SJ/JU
 Sample : J5184-07DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC26DL

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:34:51 PM

Quant Time: Oct 11 18:09:12 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 : Thu Oct 11 14:16:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2647	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	13404	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	10216	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	28284m	0.40	ng/ul	0.00
16) Chrysene-d12	21.49	240	29851m	0.40	ng/ul	0.00
20) Perylene-d12	24.05	264	27919m	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	1184	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.33	212	4372m	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.74	128	5235	0.160	ng/ul	95
5) 2-Methylnaphthalene	12.36	142	4844	0.205	ng/ul	98
7) Acenaphthylene	14.26	152	1597	0.041	ng/ul#	71
8) Acenaphthene	14.61	153	2308	0.077	ng/ul	95
9) Fluorene	15.60	166	3119	0.080	ng/ul#	95
12) Phenanthrene	17.35	178	53152	0.794	ng/ul	97
13) Anthracene	17.43	178	11550m	0.186	ng/ul	
15) Fluoranthene	19.36	202	123311m	1.434	ng/ul	
17) Pyrene	19.72	202	145826	1.640	ng/ul	99
18) Benzo(a)anthracene	21.47	228	122058m	1.352	ng/ul	
19) Chrysene	21.52	228	389268	4.800	ng/ul	96
21) Benzo(b)fluoranthene	23.26	252	286234m	3.836	ng/ul	
22) Benzo(k)fluoranthene	23.31	252	82988m	1.039	ng/ul	
23) Benzo(a)pyrene	23.93	252	95555m	1.304	ng/ul	
24) Indeno(1,2,3-cd)pyrene	26.74	276	138711m	1.668	ng/ul	
25) Dibenzo(a,h)anthracene	26.76	278	34607m	0.503	ng/ul	
26) Benzo(g,h,i)perylene	27.57	276	178147m	2.458	ng/ul	

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

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