

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE101118\
 Data File : BE097872.D
 Acq On : 12 Oct 2018 00:14
 Operator : SJ/JU
 Sample : J5183-02DL 5X
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC17DL

Manual Integrations
 APPROVED

Sohil
 10/12/2018 6:35:03 PM

Quant Time: Oct 12 05:03:13 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 04:56:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2587	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	13290	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.55	164	8881	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.31	188	21747	0.40	ng/ul	0.00
16) Chrysene-d12	21.49	240	22754	0.40	ng/ul	0.00
20) Perylene-d12	24.06	264	23304	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.29	152	991	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.34	212	3044	0.04	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	912	0.028	ng/ul#	68
5) 2-Methylnaphthalene	12.37	142	587	0.025	ng/ul	95
12) Phenanthrene	17.35	178	8411	0.163	ng/ul	95
13) Anthracene	17.44	178	1551	0.033	ng/ul#	68
15) Fluoranthene	19.37	202	14756	0.223	ng/ul	96
17) Pyrene	19.73	202	20806	0.307	ng/ul	97
18) Benzo(a)anthracene	21.48	228	12854	0.187	ng/ul#	74
19) Chrysene	21.54	228	12703	0.206	ng/ul#	78
21) Benzo(b)fluoranthene	23.27	252	16976m	0.273	ng/ul	
22) Benzo(k)fluoranthene	23.32	252	5391m	0.081	ng/ul	
23) Benzo(a)pyrene	23.95	252	10314	0.169	ng/ul#	6
24) Indeno(1,2,3-cd)pyrene	26.77	276	8591	0.124	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.78	278	2301	0.040	ng/ul#	1
26) Benzo(g,h,i)perylene	27.60	276	9806	0.162	ng/ul#	50

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

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