

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE101518\
 Data File : BE097902.D
 Acq On : 15 Oct 2018 16:31
 Operator : SJ/JU
 Sample : J5184-10DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 JLC49DL

Quant Time: Oct 15 17:08:40 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 : Mon Oct 15 16:15:33 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.90	152	1950	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	9963	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	6713	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.42	188	200219	0.40	ng/ul	0.10
16) Chrysene-d12	21.63	240	356	0.40	ng/ul	0.13
20) Perylene-d12	24.28	264	47	0.40	ng/ul	0.22
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	1193	0.07	ng/ul	0.00
14) Fluoranthene-d10	19.44	212	151	0.00	ng/ul	0.10
Target Compounds						
					Qvalue	
3) Naphthalene	10.75	128	634	0.026	ng/ul#	58
5) 2-Methylnaphthalene	12.37	142	489	0.028	ng/ul	100
8) Acenaphthene	14.62	153	633	0.032	ng/ul	96
9) Fluorene	15.61	166	518	0.020	ng/ul#	61
17) Pyrene	19.74	202	42660	40.225	ng/ul	99
18) Benzo(a)anthracene	21.67	228	2292	2.129	ng/ul#	1
19) Chrysene	21.67	228	2292	2.370	ng/ul#	1
21) Benzo(b)fluoranthene	23.48	252	4255	33.874	ng/ul#	1
22) Benzo(k)fluoranthene	23.48	252	4255	31.645	ng/ul#	1
23) Benzo(a)pyrene	24.13	252	5390	43.705	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	27.09	276	58	0.414	ng/ul#	100
25) Dibenzo(a,h)anthracene	27.12	278	2275	19.661	ng/ul#	1
26) Benzo(g,h,i)perylene	27.85	276	69	0.566	ng/ul#	1

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

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