

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE101018\
 Data File : BE097830.D
 Acq On : 10 Oct 2018 16:33
 Operator : SJ/JU
 Sample : J5295-08
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 A3Z42

Quant Time: Oct 10 18:33:57 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 : Wed Oct 10 12:43:48 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.89	152	2742	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	12477	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	8080	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.30	188	24875	0.40	ng/ul	0.00
16) Chrysene-d12	21.47	240	34744	0.40	ng/ul	0.00
20) Perylene-d12	24.03	264	34910	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	6936	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.32	212	35444	0.44	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	9838	0.328	ng/ul	99
5) 2-Methylnaphthalene	12.36	142	6590	0.330	ng/ul	98
7) Acenaphthylene	14.25	152	10358	0.364	ng/ul	97
9) Fluorene	15.59	166	13197	0.462	ng/ul	98
12) Phenanthrene	17.34	178	49987	0.841	ng/ul	99
15) Fluoranthene	19.35	202	47937	0.612	ng/ul	98
17) Pyrene	19.72	202	64642	0.757	ng/ul	99
18) Benzo(a)anthracene	21.46	228	49571	0.514	ng/ul	99
19) Chrysene	21.52	228	49251	0.529	ng/ul	99
21) Benzo(b)fluoranthene	23.25	252	32932	0.365	ng/ul	98
22) Benzo(k)fluoranthene	23.29	252	58784	0.575	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.71	276	42827	0.457	ng/ul#	100
25) Dibenzo(a,h)anthracene	26.74	278	37889	0.484	ng/ul	97
26) Benzo(g,h,i)perylene	27.53	276	59816	0.676	ng/ul	99

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

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