

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040918\
 Data File : BE095963.D
 Acq On : 9 Apr 2018 11:41
 Operator : SJ/JU
 Sample : J2135-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 X2C41

Quant Time: Apr 10 03:54:36 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 10 03:47:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1927	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	5054	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2939	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	7476	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	9928	0.40	ng/ul	0.00
20) Perylene-d12	24.40	264	7972	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	3212	0.34	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	10456	0.34	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.26	128	4497	0.319	ng/ul#	91
5) 2-Methylnaphthalene	12.84	142	5459	0.571	ng/ul	100
7) Acenaphthylene	14.70	152	3886	0.255	ng/ul	98
9) Fluorene	16.00	166	6391	0.518	ng/ul#	89
11) Pentachlorophenol	17.34	266	1707	1.079	ng/ul	96
12) Phenanthrene	17.73	178	6896	0.313	ng/ul#	89
13) Anthracene	17.81	178	10772	0.491	ng/ul#	92
15) Fluoranthene	19.69	202	8181	0.273	ng/ul	85
17) Pyrene	20.04	202	12337	0.366	ng/ul#	80
18) Benzo(a)anthracene	21.76	228	17134	0.427	ng/ul#	86
19) Chrysene	21.81	228	12520	0.434	ng/ul	97
21) Benzo(b)fluoranthene	23.58	252	18614	0.722	ng/ul	83
22) Benzo(k)fluoranthene	23.64	252	8485	0.330	ng/ul#	87
23) Benzo(a)pyrene	24.28	252	8126	0.326	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.21	278	17275	0.775	ng/ul#	86
26) Benzo(g,h,i)perylene	28.06	276	7625	0.330	ng/ul	94

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

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