

Data Path : Z:\HPCHEM1\BNA_E\Data\BE042318\
 Data File : BE096109.D
 Acq On : 24 Apr 2018 5:36
 Operator : SJ/JU
 Sample : J2434-18DL 10X
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 24 17:10:44 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 24 06:59:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	2147	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	6170	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3411	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	8882	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	7511	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	183	0.02	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	496	0.02	ng/ul	0.00

Target Compounds						Qvalue
7) Acenaphthylene	14.70	152	374	0.021	ng/ul#	71
8) Acenaphthene	15.03	153	306	0.024	ng/ul#	91
12) Phenanthrene	17.72	178	4373	0.198	ng/ul#	89
13) Anthracene	17.80	178	645	0.029	ng/ul#	85
15) Fluoranthene	19.69	202	24279	0.810	ng/ul	83
17) Pyrene	20.04	202	28125	0.932	ng/ul#	79
18) Benzo(a)anthracene	21.75	228	23652	0.659	ng/ul#	86
19) Chrysene	21.81	228	23810	0.922	ng/ul	98
21) Benzo(b)fluoranthene	23.57	252	57075	2.348	ng/ul	85
22) Benzo(k)fluoranthene	23.57	252	57075	2.357	ng/ul#	87
23) Benzo(a)pyrene	24.26	252	31831	1.357	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.18	276	23254	0.892	ng/ul#	74
25) Dibenzo(a,h)anthracene	27.18	278	5791	0.276	ng/ul	98
26) Benzo(g,h,i)perylene	28.05	276	24545	1.127	ng/ul#	93

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

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