

Data Path : Z:\HPCHEM1\BNA_E\Data\BE042318\
 Data File : BE096102.D
 Acq On : 24 Apr 2018 1:25
 Operator : SJ/JU
 Sample : J2434-06
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Apr 24 14:57:36 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 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	2209	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	6181	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3784	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.77	188	323314	0.40	ng/ul	0.10
16) Chrysene-d12	21.89	240	374	0.40	ng/ul	0.12
20) Perylene-d12	24.40	264	11370	0.40	ng/ul	0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	2154	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.82	212	163	0.00	ng/ul	0.16
Target Compounds				Qvalue		
3) Naphthalene	11.25	128	29440	1.706	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	23012	1.966	ng/ul	100
7) Acenaphthylene	14.70	152	49999	2.549	ng/ul	96
8) Acenaphthene	15.03	153	5912	0.420	ng/ul#	90
9) Fluorene	15.99	166	8877	0.558	ng/ul	88
12) Phenanthrene	17.81	178	113613	0.119	ng/ul#	89
15) Fluoranthene	19.83	202	27322	0.021	ng/ul	84
17) Pyrene	20.05	202	1044012	821.497	ng/ul#	75
18) Benzo(a)anthracene	21.82	228	690567	457.041	ng/ul#	83
19) Chrysene	21.82	228	689475	634.018	ng/ul	95
21) Benzo(b)fluoranthene	23.60	252	1637857	44.517	ng/ul	83
22) Benzo(k)fluoranthene	23.60	252	1637857	44.686	ng/ul#	84
23) Benzo(a)pyrene	24.29	252	603300	16.993	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.22	276	450076	11.407	ng/ul#	73
25) Dibenzo(a,h)anthracene	27.34	278	38118	1.200	ng/ul	98
26) Benzo(g,h,i)perylene	28.09	276	414836	12.587	ng/ul#	90

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

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