

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006566.D
 Acq On : 25 Jun 2019 11:38
 Operator : HP/JU
 Sample : K3362-22
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41R6

Quant Time: Jun 25 12:25:15 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 25 07:15:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	4635	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	17593	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9037	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	17352	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9628	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	10136	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	7354	0.27	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	13810	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	11423	0.235	ng/ul	98
5) 2-Methylnaphthalene	12.08	142	7134	0.211	ng/ul	99
7) Acenaphthylene	14.00	152	13876	0.276	ng/ul	99
9) Fluorene	15.34	166	11003	0.270	ng/ul	97
12) Phenanthrene	17.08	178	33505	0.602	ng/ul	99
15) Fluoranthene	19.11	202	21806	0.355	ng/ul	96
17) Pyrene	19.48	202	23434	0.462	ng/ul	100
18) Benzo(a)anthracene	21.24	228	11198	0.263	ng/ul	99
19) Chrysene	21.30	228	14367	0.359	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	10309	0.244	ng/ul	94
22) Benzo(k)fluoranthene	22.87	252	16566	0.402	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.74	276	14218	0.334	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.74	278	11885	0.349	ng/ul	98
26) Benzo(g,h,i)perylene	26.41	276	18744	0.525	ng/ul	96

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

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