

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071919\
 Data File : BN006921.D
 Acq On : 19 Jul 2019 13:15
 Operator : HP/JU
 Sample : K3362-22
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41R6

Quant Time: Jul 19 23:32:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 19 23:29:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	3527	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.37	136	12073	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	6764	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	14514	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	12169	0.40	ng/ul	0.00
20) Perylene-d12	23.47	264	13323	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	4832	0.26	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	12827	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	7706	0.222	ng/ul	95
5) 2-Methylnaphthalene	12.06	142	4658	0.199	ng/ul	99
7) Acenaphthylene	13.96	152	8637	0.247	ng/ul	99
9) Fluorene	15.31	166	7524	0.255	ng/ul	99
12) Phenanthrene	17.05	178	25961	0.556	ng/ul	100
15) Fluoranthene	19.08	202	19718	0.342	ng/ul	99
17) Pyrene	19.45	202	22057	0.397	ng/ul	98
18) Benzo(a)anthracene	21.22	228	11955	0.239	ng/ul	99
19) Chrysene	21.27	228	17717	0.318	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	11862	0.217	ng/ul	97
22) Benzo(k)fluoranthene	22.84	252	21180	0.371	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.69	276	17116	0.316	ng/ul	95
25) Dibenzo(a,h)anthracene	25.69	278	14214	0.330	ng/ul	98
26) Benzo(g,h,i)perylene	26.37	276	22961	0.507	ng/ul	96

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

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