

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006896.D
 Acq On : 18 Jul 2019 17:52
 Operator : HP/JU
 Sample : K3362-22
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41R6

Quant Time: Jul 19 03:52:32 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 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	1819	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6931	0.40	ng/ul	0.01
6) Acenaphthene-d10	14.24	164	3907	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	8342	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7349	0.40	ng/ul	-0.01
20) Perylene-d12	23.45	264	8087	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	3373	0.32	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	9287	0.38	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	6073	0.305	ng/ul#	94
5) 2-Methylnaphthalene	12.06	142	3356	0.250	ng/ul	99
7) Acenaphthylene	13.96	152	6393	0.317	ng/ul	98
9) Fluorene	15.31	166	5434	0.319	ng/ul	99
12) Phenanthrene	17.05	178	18837	0.702	ng/ul	99
15) Fluoranthene	19.08	202	14184	0.428	ng/ul	97
17) Pyrene	19.44	202	15919	0.474	ng/ul	99
18) Benzo(a)anthracene	21.21	228	8874	0.294	ng/ul	100
19) Chrysene	21.27	228	12401	0.368	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	8465	0.255	ng/ul	93
22) Benzo(k)fluoranthene	22.83	252	15171	0.437	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.68	276	12407	0.378	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.67	278	10451	0.400	ng/ul	98
26) Benzo(g,h,i)perylene	26.35	276	16480	0.600	ng/ul	97

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

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