

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN071819\
 Data File : BN006911.D
 Acq On : 19 Jul 2019 03:42
 Operator : HP/JU
 Sample : K3764-15
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A52F9

Manual Integrations
 APPROVED

mohammad
 7/19/2019 4:00:06 PM

Quant Time: Jul 19 05:45:51 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 03:57:16 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	3455	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.35	136	12530	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.23	164	6419	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.00	188	13804	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	11183	0.40	ng/ul	-0.02
20) Perylene-d12	23.45	264	14867	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	1484	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	4316	0.11	ng/ul	-0.01
Target Compounds				Ovalue		
12) Phenanthrene	17.04	178	2582	0.058	ng/ul	95
15) Fluoranthene	19.08	202	7039	0.128	ng/ul	99
17) Pyrene	19.43	202	6159m	0.121	ng/ul	
18) Benzo(a)anthracene	21.22	228	2162m	0.047	ng/ul	
19) Chrysene	21.27	228	3238	0.063	ng/ul	94
21) Benzo(b)fluoranthene	22.79	252	4756m	0.078	ng/ul	
22) Benzo(k)fluoranthene	22.83	252	1454m	0.023	ng/ul	
23) Benzo(a)pyrene	23.34	252	2715	0.045	ng/ul#	71
24) Indeno(1,2,3-cd)pyrene	25.70	276	2172	0.036	ng/ul#	87
26) Benzo(g,h,i)perylene	26.37	276	2314	0.046	ng/ul#	86

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

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