

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080719\
 Data File : BN006997.D
 Acq On : 07 Aug 2019 20:57
 Operator : HP/JU
 Sample : K3893-15DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP19DL

Manual Integrations
 APPROVED

mohammad
 8/9/2019 6:15:11 AM

Quant Time: Aug 08 05:24:09 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN080719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 07 17:30:08 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	7344	0.40	ng/ul	0.00
2) Naphthalene-d8	10.23	136	29517	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.12	164	18825	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	47819	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	40019	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	35525	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.83	152	2189	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.90	212	6447	0.04	ng/ul	0.00
Target Compounds				Qvalue		
7) Acenaphthylene	13.83	152	6965	0.064	ng/ul	97
9) Fluorene	15.17	166	3713	0.041	ng/ul	99
12) Phenanthrene	16.90	178	66153	0.434	ng/ul	99
13) Anthracene	16.99	178	10576	0.068	ng/ul	97
15) Fluoranthene	18.93	202	135275	0.714	ng/ul	98
17) Pyrene	19.29	202	99920	0.559	ng/ul	99
18) Benzo(a)anthracene	21.06	228	54825	0.312	ng/ul	96
19) Chrysene	21.11	228	54910	0.334	ng/ul	97
21) Benzo(b)fluoranthene	22.58	252	61684m	0.428	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	23998m	0.172	ng/ul	
23) Benzo(a)pyrene	23.11	252	37773	0.279	ng/ul#	96
24) Indeno(1,2,3-cd)pyrene	25.30	276	24504	0.168	ng/ul#	94
25) Dibenzo(a,h)anthracene	25.31	278	6474	0.055	ng/ul#	74
26) Benzo(g,h,i)perylene	25.94	276	20666	0.169	ng/ul	97

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

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