

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080919\
 Data File : BN007041.D
 Acq On : 09 Aug 2019 16:10
 Operator : HP/JU
 Sample : K3962-15
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP08

Manual Integrations
 APPROVED

mohammad
 8/14/2019 2:49:44 AM

Quant Time: Aug 09 17:29:11 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	7.46	152	9262	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.22	136	35205	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.11	164	20995	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.86	188	53791m	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	46021m	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	39496m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.82	152	16154	0.26	ng/ul	-0.01
14) Fluoranthene-d10	18.90	212	48659m	0.28	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.27	128	3247	0.032	ng/ul#	93
5) 2-Methylnaphthalene	11.90	142	1771	0.023	ng/ul	99
7) Acenaphthylene	13.82	152	6576	0.054	ng/ul	97
8) Acenaphthene	14.17	153	2014	0.024	ng/ul	98
9) Fluorene	15.17	166	3355	0.033	ng/ul#	96
12) Phenanthrene	16.90	178	77745m	0.454	ng/ul	
13) Anthracene	16.99	178	19359m	0.110	ng/ul	
15) Fluoranthene	18.93	202	160532m	0.753	ng/ul	
17) Pyrene	19.30	202	171048	0.833	ng/ul	98
18) Benzo(a)anthracene	21.06	228	103372m	0.511	ng/ul	
19) Chrysene	21.12	228	102010	0.539	ng/ul	98
21) Benzo(b)fluoranthene	22.59	252	98823m	0.617	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	41313m	0.266	ng/ul	
23) Benzo(a)pyrene	23.12	252	77179	0.512	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.31	276	44374	0.273	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.33	278	14626	0.113	ng/ul#	79
26) Benzo(g,h,i)perylene	25.94	276	43759	0.321	ng/ul	98

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

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