

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM081219\
 Data File : BM022023.D
 Acq On : 12 Aug 2019 09:47
 Operator : HP/JU
 Sample : K4028-03DL 5X
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BEP69DL

Manual Integrations
 APPROVED

Sohil
 8/13/2019 9:56:11 AM

Quant Time: Aug 12 14:14:52 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 09 01:29:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	647	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.40	136	2742	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.27	164	1485	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.02	188	3341	0.40	ng/ul	-0.02
16) Chrysene-d12	21.22	240	3422	0.40	ng/ul	-0.02
20) Perylene-d12	23.43	264	3652m	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	150	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	471	0.04	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	273	0.035	ng/ul#	45
5) 2-Methylnaphthalene	12.09	142	169	0.033	ng/ul	93
7) Acenaphthylene	13.99	152	1135	0.164	ng/ul#	81
8) Acenaphthene	14.33	153	735	0.130	ng/ul	95
9) Fluorene	15.33	166	1013	0.161	ng/ul#	98
12) Phenanthrene	17.06	178	21090	2.169	ng/ul	97
13) Anthracene	17.16	178	5577	0.673	ng/ul	98
15) Fluoranthene	19.08	202	46475	3.569	ng/ul	97
17) Pyrene	19.45	202	52984	3.768	ng/ul	99
18) Benzo(a)anthracene	21.20	228	31972	2.612	ng/ul	99
19) Chrysene	21.26	228	33471	2.126	ng/ul	99
21) Benzo(b)fluoranthene	22.77	252	33230m	2.698	ng/ul	
22) Benzo(k)fluoranthene	22.81	252	13078m	0.938	ng/ul	
23) Benzo(a)pyrene	23.33	252	27603	2.204	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	25.64	276	15638	1.046	ng/ul#	95
25) Dibenzo(a,h)anthracene	25.64	278	4700	0.389	ng/ul	95
26) Benzo(a,h,i)perylene	26.32	276	16251	1.207	ng/ul	98

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

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