

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN060619\
 Data File : BN006174.D
 Acq On : 07 Jun 2019 19:19
 Operator : JU/SJ
 Sample : K3201-10DL 5X
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC1DL

Manual Integrations
 APPROVED

mohammad
 6/11/2019 5:08:48 PM

Quant Time: Jun 08 03:15:39 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 07 01:31:14 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.65	152	5142	0.40	ng/ul	0.00
2) Naphthalene-d8	10.43	136	19597	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.30	164	9679	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	19826	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	16703	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	21504	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.03	152	1567	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	3278	0.06	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.48	128	1791	0.033	ng/ul#	85
5) 2-Methylnaphthalene	12.10	142	804	0.021	ng/ul	99
7) Acenaphthylene	14.01	152	8231	0.153	ng/ul	98
8) Acenaphthene	14.36	153	2309	0.062	ng/ul	99
9) Fluorene	15.35	166	2626	0.060	ng/ul#	92
12) Phenanthrene	17.09	178	67255	1.057	ng/ul	100
13) Anthracene	17.18	178	14983	0.247	ng/ul	98
15) Fluoranthene	19.12	202	175105	2.492	ng/ul	99
17) Pyrene	19.49	202	169196	1.923	ng/ul	99
18) Benzo(a)anthracene	21.25	228	86783	1.175	ng/ul	100
19) Chrysene	21.31	228	81660	1.177	ng/ul	98
21) Benzo(b)fluoranthene	22.84	252	101325m	1.130	ng/ul	
22) Benzo(k)fluoranthene	22.88	252	43812m	0.502	ng/ul	
23) Benzo(a)pyrene	23.40	252	71265	0.853	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.73	276	48889	0.541	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.73	278	12254	0.170	ng/ul	99
26) Benzo(a,h,i)perylene	26.42	276	37845	0.500	ng/ul	96

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

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