

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062219\
 Data File : BN006498.D
 Acq On : 22 Jun 2019 20:24
 Operator : HP/JU
 Sample : K3335-08DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4235DL

Manual Integrations
 APPROVED

mohammad
 6/25/2019 8:44:39 AM

Quant Time: Jun 24 10:39:23 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 : Mon Jun 17 11:13:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	4068	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.41	136	16067	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.28	164	8730	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	22987	0.40	ng/ul	-0.01
16) Chrysene-d12	21.25	240	13559	0.40	ng/ul	-0.02
20) Perylene-d12	23.48	264	13080	0.40	ng/ul	-0.03
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	952	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2137	0.03	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	2250	0.051	ng/ul#	88
5) 2-Methylnaphthalene	12.09	142	814	0.026	ng/ul	95
7) Acenaphthylene	14.00	152	14311	0.295	ng/ul	99
8) Acenaphthene	14.34	153	3825	0.114	ng/ul	98
9) Fluorene	15.34	166	7308	0.186	ng/ul	97
12) Phenanthrene	17.08	178	112817	1.529	ng/ul	100
13) Anthracene	17.17	178	26358	0.374	ng/ul	99
15) Fluoranthene	19.11	202	241725	2.967	ng/ul	99
17) Pyrene	19.47	202	213233	2.986	ng/ul	99
18) Benzo(a)anthracene	21.24	228	86636	1.445	ng/ul	100
19) Chrysene	21.29	228	81415	1.446	ng/ul	98
21) Benzo(b)fluoranthene	22.81	252	92798m	1.701	ng/ul	
22) Benzo(k)fluoranthene	22.85	252	36528m	0.688	ng/ul	
23) Benzo(a)pyrene	23.38	252	61899	1.219	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.71	276	37059	0.674	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.71	278	9128	0.208	ng/ul	97
26) Benzo(g,h,i)perylene	26.40	276	30143	0.655	ng/ul	95

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

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