

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102219\
 Data File : BN008357.D
 Acq On : 23 Oct 2019 19:44
 Operator : JU
 Sample : K5223-10DL 2X
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPE2DL

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:06:51 AM

Quant Time: Oct 24 01:57:51 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1935	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	9333	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5983	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12857	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	10701	0.40	ng/ul	-0.01
20) Perylene-d12	23.34	264	8627	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	1393	0.09	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	4116	0.09	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	1025	0.039	ng/ul#	85
5) 2-Methylnaphthalene	12.02	142	590	0.032	ng/ul	99
7) Acenaphthylene	13.93	152	703	0.025	ng/ul#	75
8) Acenaphthene	14.28	153	2634	0.119	ng/ul	98
9) Fluorene	15.27	166	2766	0.110	ng/ul	98
12) Phenanthrene	17.00	178	46864	1.245	ng/ul	99
13) Anthracene	17.09	178	9432	0.285	ng/ul	98
15) Fluoranthene	19.03	202	59995	1.207	ng/ul	96
17) Pyrene	19.39	202	54221	1.273	ng/ul	99
18) Benzo(a)anthracene	21.15	228	26916	0.692	ng/ul	97
19) Chrysene	21.20	228	26416	0.553	ng/ul	97
21) Benzo(b)fluoranthene	22.70	252	25539m	0.909	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	9246m	0.291	ng/ul	
23) Benzo(a)pyrene	23.25	252	18699	0.633	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.49	276	10184	0.293	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.49	278	3305	0.120	ng/ul#	79
26) Benzo(g,h,i)perylene	26.14	276	9087	0.285	ng/ul	95

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

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