

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102219\
 Data File : BN008359.D
 Acq On : 23 Oct 2019 20:55
 Operator : JU
 Sample : K5223-17DL 2X
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPA5DL

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:06:57 AM

Quant Time: Oct 24 02:03:02 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	2080	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	10028	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	6359	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	13781	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	11905	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	10063	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	1779	0.11	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	5239	0.11	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	640	0.022	ng/ul#	73
7) Acenaphthylene	13.93	152	936	0.031	ng/ul#	83
8) Acenaphthene	14.27	153	744	0.032	ng/ul	94
9) Fluorene	15.27	166	799	0.030	ng/ul#	95
12) Phenanthrene	17.00	178	15494	0.384	ng/ul	100
13) Anthracene	17.09	178	3001	0.085	ng/ul#	94
15) Fluoranthene	19.03	202	35733	0.671	ng/ul	96
17) Pyrene	19.40	202	41901	0.884	ng/ul	99
18) Benzo(a)anthracene	21.15	228	20591	0.476	ng/ul	97
19) Chrysene	21.21	228	23611	0.445	ng/ul	98
21) Benzo(b)fluoranthene	22.70	252	21588m	0.659	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	8172m	0.220	ng/ul	
23) Benzo(a)pyrene	23.25	252	16156	0.469	ng/ul#	90
24) Indeno(1,2,3-cd)pyrene	25.49	276	8638	0.213	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.49	278	2100	0.065	ng/ul#	71
26) Benzo(g,h,i)perylene	26.14	276	6957	0.187	ng/ul#	94

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

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