

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN102219\
 Data File : BN008355.D
 Acq On : 23 Oct 2019 18:32
 Operator : JU
 Sample : K5223-05DL 2X
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPD9DL

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:06:46 AM

Quant Time: Oct 24 01:52:13 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	1902	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.34	136	9250	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5880	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12673	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	10584	0.40	ng/ul	0.00
20) Perylene-d12	23.34	264	8589	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.95	152	1572	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.00	212	4482	0.10	ng/ul	-0.01
Target Compounds				Qvalue		
3) Naphthalene	10.39	128	602	0.023	ng/ul#	70
7) Acenaphthylene	13.93	152	1324	0.047	ng/ul#	60
8) Acenaphthene	14.28	153	491	0.023	ng/ul	95
9) Fluorene	15.27	166	643	0.026	ng/ul#	91
12) Phenanthrene	17.00	178	14920	0.402	ng/ul	100
13) Anthracene	17.10	178	2752	0.084	ng/ul	94
15) Fluoranthene	19.03	202	40080	0.818	ng/ul	98
17) Pyrene	19.39	202	37668	0.894	ng/ul	98
18) Benzo(a)anthracene	21.16	228	17492	0.455	ng/ul	97
19) Chrysene	21.21	228	21412	0.453	ng/ul	98
21) Benzo(b)fluoranthene	22.70	252	22683m	0.811	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	7585m	0.239	ng/ul	
23) Benzo(a)pyrene	23.25	252	14510	0.493	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.49	276	9012	0.260	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.50	278	1967	0.072	ng/ul#	71
26) Benzo(g,h,i)perylene	26.15	276	6929	0.218	ng/ul	95

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

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