

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101819\
 Data File : BN008251.D
 Acq On : 18 Oct 2019 17:24
 Operator : JU
 Sample : K5178-05DL 5X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPE8DL

Manual Integrations
 APPROVED

mohammad
 10/22/2019 1:40:32 AM

Quant Time: Oct 18 18:08:12 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 : Fri Oct 18 17:22:25 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	3675	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	17696	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.23	164	11193	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.98	188	23515	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	22079	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	19051	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	778	0.03	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	2406	0.03	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.95	152	1336	0.025	ng/ul#	87
12) Phenanthrene	17.02	178	10102	0.147	ng/ul	99
13) Anthracene	17.11	178	1950	0.032	ng/ul#	90
15) Fluoranthene	19.05	202	23038	0.253	ng/ul	99
17) Pyrene	19.41	202	24902	0.283	ng/ul	99
18) Benzo(a)anthracene	21.17	228	13200	0.164	ng/ul	95
19) Chrysene	21.22	228	13791	0.140	ng/ul	97
21) Benzo(b)fluoranthene	22.72	252	15661m	0.253	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	6371m	0.091	ng/ul	
23) Benzo(a)pyrene	23.26	252	11064	0.170	ng/ul#	89
24) Indeno(1,2,3-cd)pyrene	25.52	276	6507	0.085	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.53	278	1547	0.025	ng/ul#	45
26) Benzo(g,h,i)perylene	26.17	276	6382	0.091	ng/ul#	91

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

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