

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008284.D
 Acq On : 21 Oct 2019 15:48
 Operator : JU
 Sample : K5199-13
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEP80

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:27:21 PM

Quant Time: Oct 21 18:38:39 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 : Mon Oct 21 16:47:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1412	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	6789	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	4258	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	9581m	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	9127	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	11559	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	3582	0.31	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	10695	0.32	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.01	178	1994	0.071	ng/ul#	93
15) Fluoranthene	19.04	202	4177	0.113	ng/ul	95
17) Pyrene	19.40	202	3477	0.096	ng/ul#	88
18) Benzo(a)anthracene	21.16	228	1908	0.057	ng/ul#	88
19) Chrysene	21.22	228	2108	0.052	ng/ul#	91
21) Benzo(b)fluoranthene	22.71	252	2879m	0.077	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	1123m	0.026	ng/ul	
23) Benzo(a)pyrene	23.26	252	1731	0.044	ng/ul#	20
24) Indeno(1,2,3-cd)pyrene	25.51	276	1042	0.022	ng/ul	92
26) Benzo(g,h,i)perylene	26.17	276	865	0.020	ng/ul#	51

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

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