

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008292.D
 Acq On : 21 Oct 2019 20:38
 Operator : JU
 Sample : K5178-09DL 10X
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPF2DL

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:27:48 PM

Quant Time: Oct 22 00:30:25 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 23:58:15 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	1824	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	8495	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5482	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	11170	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	10629	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	9016	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	362	0.03	ng/ul	0.01
14) Fluoranthene-d10	19.01	212	1137	0.03	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.01	178	2293	0.070	ng/ul#	93
13) Anthracene	17.11	178	602	0.021	ng/ul#	64
15) Fluoranthene	19.04	202	4268	0.099	ng/ul	95
17) Pyrene	19.40	202	4438	0.105	ng/ul#	90
18) Benzo(a)anthracene	21.16	228	2957	0.077	ng/ul#	89
19) Chrysene	21.22	228	2439	0.051	ng/ul#	92
21) Benzo(b)fluoranthene	22.71	252	3583m	0.122	ng/ul	
22) Benzo(k)fluoranthene	22.74	252	1274m	0.038	ng/ul	
23) Benzo(a)pyrene	23.26	252	2025	0.066	ng/ul#	38
24) Indeno(1,2,3-cd)pyrene	25.51	276	1182	0.032	ng/ul#	96
26) Benzo(g,h,i)perylene	26.16	276	1259	0.038	ng/ul#	65

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

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