

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102119\
 Data File : BN008290.D
 Acq On : 21 Oct 2019 19:27
 Operator : JU
 Sample : K5178-21DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPJ3DL

Manual Integrations
 APPROVED

mohammad
 10/22/2019 3:27:43 PM

Quant Time: Oct 22 00:24:16 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	1965	0.40	ng/ul	0.00
2) Naphthalene-d8	10.35	136	9314	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	5968	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.97	188	12730	0.40	ng/ul	0.00
16) Chrysene-d12	21.18	240	13094	0.40	ng/ul	0.00
20) Perylene-d12	23.35	264	11282	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	840	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.01	212	2921	0.07	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.40	128	1082	0.041	ng/ul#	83
5) 2-Methylnaphthalene	12.03	142	590	0.032	ng/ul	95
7) Acenaphthylene	13.94	152	3596	0.127	ng/ul	96
8) Acenaphthene	14.29	153	1460	0.066	ng/ul	98
9) Fluorene	15.28	166	1964	0.078	ng/ul#	97
12) Phenanthrene	17.01	178	44099	1.183	ng/ul	99
13) Anthracene	17.10	178	8024	0.245	ng/ul	98
15) Fluoranthene	19.04	202	119983	2.437	ng/ul	97
17) Pyrene	19.40	202	120343	2.308	ng/ul	100
18) Benzo(a)anthracene	21.16	228	63896	1.342	ng/ul	98
19) Chrysene	21.22	228	74835	1.281	ng/ul	99
21) Benzo(b)fluoranthene	22.71	252	81089m	2.208	ng/ul	
22) Benzo(k)fluoranthene	22.75	252	27744m	0.667	ng/ul	
23) Benzo(a)pyrene	23.26	252	55933	1.447	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.51	276	31745	0.697	ng/ul	99
25) Dibenzo(a,h)anthracene	25.51	278	7503	0.209	ng/ul	92
26) Benzo(a,h,i)perylene	26.16	276	30071	0.720	ng/ul	98

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

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