

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009245.D
 Acq On : 29 Dec 2019 01:16
 Operator : JU
 Sample : K6278-20DL 5X
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C84DL

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:33:31 PM

Quant Time: Dec 29 05:56:16 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Dec 29 04:38:59 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2599	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	9433	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4997	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	10102	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	8027	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	8159	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	651	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	1286	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	1151	0.040	ng/ul#	86
5) 2-Methylnaphthalene	11.94	142	838	0.039	ng/ul	99
7) Acenaphthylene	13.86	152	1529	0.058	ng/ul#	90
9) Fluorene	15.19	166	633	0.026	ng/ul#	91
12) Phenanthrene	16.93	178	13575	0.386	ng/ul	99
13) Anthracene	17.02	178	2547	0.081	ng/ul#	93
15) Fluoranthene	18.95	202	37566	0.935	ng/ul	100
17) Pyrene	19.32	202	31343	0.765	ng/ul	99
18) Benzo(a)anthracene	21.07	228	14487	0.406	ng/ul	97
19) Chrysene	21.12	228	18029	0.503	ng/ul	97
21) Benzo(b)fluoranthene	22.59	252	23720m	0.623	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	8150m	0.212	ng/ul	
23) Benzo(a)pyrene	23.12	252	14305	0.428	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.30	276	10850	0.298	ng/ul#	89
25) Dibenzo(a,h)anthracene	25.31	278	2538	0.088	ng/ul#	67
26) Benzo(a,h,i)perylene	25.94	276	10975	0.358	ng/ul	98

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

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