

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN123019\
 Data File : BN009270.D
 Acq On : 30 Dec 2019 21:18
 Operator : JU
 Sample : K6322-11DL 5X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C95DL

Manual Integrations
 APPROVED

mohammad
 12/31/2019 12:12:46 PM

Quant Time: Dec 31 01:58:47 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 : Tue Dec 31 00:59:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	2428	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	8550	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.13	164	4511	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.88	188	8889	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	6975	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7113	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	534	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	1019	0.04	ng/ul	0.00
Target Compounds				Ovalue		
7) Acenaphthylene	13.85	152	1788	0.075	ng/ul#	93
9) Fluorene	15.19	166	1084	0.049	ng/ul#	96
12) Phenanthrene	16.92	178	17587	0.569	ng/ul	100
13) Anthracene	17.02	178	3332	0.120	ng/ul	97
15) Fluoranthene	18.95	202	32393	0.916	ng/ul	100
17) Pyrene	19.31	202	27585	0.775	ng/ul	99
18) Benzo(a)anthracene	21.07	228	10789	0.348	ng/ul	98
19) Chrysene	21.12	228	10578	0.340	ng/ul	97
21) Benzo(b)fluoranthene	22.59	252	12999m	0.391	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	4918m	0.147	ng/ul	
23) Benzo(a)pyrene	23.12	252	9508	0.327	ng/ul	96
24) Indeno(1,2,3-cd)pyrene	25.30	276	6341	0.200	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.31	278	1405	0.056	ng/ul#	62
26) Benzo(g,h,i)perylene	25.93	276	6441	0.241	ng/ul	98

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

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