

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009248.D
 Acq On : 29 Dec 2019 03:04
 Operator : JU
 Sample : K6278-15
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C68

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:33:41 PM

Quant Time: Dec 29 06:04:00 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	2508	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	9056	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4785	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	9663m	0.40	ng/ul	0.00
16) Chrysene-d12	21.08	240	7666	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7843	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	3250	0.20	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	6201	0.22	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	4669	0.170	ng/ul	96
5) 2-Methylnaphthalene	11.93	142	2496	0.120	ng/ul	100
7) Acenaphthylene	13.85	152	4343	0.173	ng/ul	95
8) Acenaphthene	14.20	153	651	0.033	ng/ul	95
9) Fluorene	15.19	166	1637	0.070	ng/ul#	96
12) Phenanthrene	16.93	178	19501	0.580	ng/ul	99
13) Anthracene	17.02	178	5423	0.180	ng/ul	97
15) Fluoranthene	18.95	202	39676	1.032	ng/ul	99
17) Pyrene	19.31	202	32110	0.821	ng/ul	98
18) Benzo(a)anthracene	21.07	228	17712	0.520	ng/ul	94
19) Chrysene	21.12	228	16856m	0.493	ng/ul	
21) Benzo(b)fluoranthene	22.58	252	22010m	0.601	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	8144m	0.221	ng/ul	
23) Benzo(a)pyrene	23.12	252	14352	0.447	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.30	276	10100	0.289	ng/ul#	88
25) Dibenzo(a,h)anthracene	25.30	278	2621	0.094	ng/ul#	63
26) Benzo(a,h,i)perylene	25.93	276	10094	0.343	ng/ul	98

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

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