

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122819\
 Data File : BN009243.D
 Acq On : 29 Dec 2019 00:04
 Operator : JU
 Sample : K6278-17DL 5X
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0C70DL

Manual Integrations
 APPROVED

mohammad
 12/29/2019 5:33:23 PM

Quant Time: Dec 29 05:51:53 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	2478	0.40	ng/ul	0.00
2) Naphthalene-d8	10.26	136	8979	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.14	164	4761	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.89	188	9612	0.40	ng/ul	0.00
16) Chrysene-d12	21.09	240	7670	0.40	ng/ul	0.00
20) Perylene-d12	23.21	264	7948	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.86	152	578	0.04	ng/ul	0.00
14) Fluoranthene-d10	18.92	212	1169	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.31	128	824	0.030	ng/ul#	81
5) 2-Methylnaphthalene	11.94	142	564	0.027	ng/ul	96
12) Phenanthrene	16.93	178	2513	0.075	ng/ul#	93
15) Fluoranthene	18.95	202	3946	0.103	ng/ul	95
17) Pyrene	19.32	202	3659	0.094	ng/ul	97
18) Benzo(a)anthracene	21.07	228	1839	0.054	ng/ul#	86
19) Chrysene	21.12	228	2485	0.073	ng/ul#	89
21) Benzo(b)fluoranthene	22.59	252	3360m	0.091	ng/ul	
22) Benzo(k)fluoranthene	22.62	252	1478m	0.039	ng/ul	
23) Benzo(a)pyrene	23.12	252	1941	0.060	ng/ul#	41
24) Indeno(1,2,3-cd)pyrene	25.30	276	1577	0.045	ng/ul#	90
26) Benzo(g,h,i)perylene	25.94	276	1749	0.059	ng/ul#	72

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

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