

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012821\
 Data File : BN013526.D
 Acq On : 28 Jan 2021 13:57
 Operator : CG/JU
 Sample : M1203-09
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2Y85

Quant Time: Jan 28 14:38:18 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN012521.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 28 11:56:28 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	4024	0.40	ng/ul	0.00
4) Naphthalene-d8	10.33	136	15252	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.20	164	8062	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.95	188	17666	0.40	ng/ul	0.00
17) Chrysene-d12	21.15	240	15733	0.40	ng/ul	0.00
23) Perylene-d12	23.32	264	13555	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.17	96	1756	0.43	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.92	152	6359	0.28	ng/ul	0.00
18) Fluoranthene-d10	18.99	212	16132	0.34	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.38	128	10478	0.244	ng/ul	99
7) 2-Methylnaphthalene	12.00	142	13757	0.477	ng/ul	100
10) Acenaphthylene	13.92	152	8230	0.243	ng/ul	99
11) Acenaphthene	14.26	153	9424	0.324	ng/ul	99
12) Fluorene	15.25	166	15694	0.486	ng/ul	98
14) Pentachlorophenol	16.60	266	2213	1.037	ng/ul	99
15) Phenanthrene	16.99	178	20903	0.368	ng/ul	99
16) Anthracene	17.08	178	23016	0.499	ng/ul	99
19) Fluoranthene	19.01	202	23740	0.381	ng/ul	98
20) Pyrene	19.38	202	23639	0.378	ng/ul	99
21) Benzo(a)anthracene	21.14	228	27520	0.555	ng/ul	99
22) Chrysene	21.19	228	27411	0.446	ng/ul	99
24) Benzo(b)fluoranthene	22.67	252	34449	0.582	ng/ul	100
25) Benzo(k)fluoranthene	22.72	252	18608	0.302	ng/ul	98
26) Benzo(a)pyrene	23.23	252	22212	0.477	ng/ul	98
28) Dibenzo(a,h)anthracene	25.49	278	36029	0.677	ng/ul	99
29) Benzo(g,h,i)perylene	26.13	276	15797	0.273	ng/ul	98

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

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