

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN012821\
 Data File : BN013527.D
 Acq On : 28 Jan 2021 14:53
 Operator : CG/JU
 Sample : M1203-04
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X2Y80

Quant Time: Jan 28 15:26:49 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	7.56	152	4279	0.40	ng/ul	0.00
4) Naphthalene-d8	10.32	136	16045	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.20	164	8462	0.40	ng/ul	0.00
13) Phenanthrene-d10	16.95	188	18063	0.40	ng/ul	0.00
17) Chrysene-d12	21.15	240	15280	0.40	ng/ul	0.00
23) Perylene-d12	23.33	264	13597	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.17	96	2647	0.61	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.92	152	8074	0.34	ng/ul	0.00
18) Fluoranthene-d10	18.99	212	17241	0.38	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.37	128	16318	0.361	ng/ul	100
7) 2-Methylnaphthalene	11.99	142	23101	0.762	ng/ul	99
10) Acenaphthylene	13.92	152	17218	0.484	ng/ul	100
11) Acenaphthene	14.26	153	34323	1.124	ng/ul	100
14) Pentachlorophenol	16.60	266	1328	0.609	ng/ul	99
15) Phenanthrene	16.99	178	29758	0.513	ng/ul	99
19) Fluoranthene	19.02	202	54849	0.906	ng/ul	100
20) Pyrene	19.38	202	43641	0.719	ng/ul	99
21) Benzo(a)anthracene	21.14	228	26119	0.543	ng/ul	100
24) Benzo(b)fluoranthene	22.68	252	45143	0.760	ng/ul	98
25) Benzo(k)fluoranthene	22.72	252	55919	0.904	ng/ul	97
26) Benzo(a)pyrene	23.23	252	25523	0.547	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.49	276	47838	0.729	ng/ul#	78
28) Dibenzo(a,h)anthracene	25.50	278	48140	0.902	ng/ul	99
29) Benzo(a,h,i)perylene	26.14	276	52535	0.906	ng/ul	99

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

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