

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN041321\
 Data File : BN014270.D
 Acq On : 13 Apr 2021 14:19
 Operator : CG/JU
 Sample : M1906-04
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3309

Quant Time: Apr 14 01:36:25 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SFAM-EPA-SIM-BN041221.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 12 17:52:31 2021
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.70	152	5732	0.40	ng/ul	0.00
4) Naphthalene-d8	10.48	136	19120	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.34	164	9699	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.08	188	18950	0.40	ng/ul	0.00
17) Chrysene-d12	21.28	240	15412	0.40	ng/ul	0.00
23) Perylene-d12	23.52	264	13023	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.23	96	4678	0.75	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.07	152	11417	0.39	ng/ul	0.00
18) Fluoranthene-d10	19.12	212	19938	0.42	ng/ul	0.00
Target Compounds				Ovalue		
5) Naphthalene	10.53	128	20690	0.366	ng/ul	100
7) 2-Methylnaphthalene	12.14	142	23343	0.647	ng/ul	100
10) Acenaphthylene	14.06	152	11759	0.259	ng/ul	99
12) Fluorene	15.39	166	22484	0.570	ng/ul	99
14) Pentachlorophenol	16.73	266	4453	0.726	ng/ul	99
15) Phenanthrene	17.12	178	22538	0.368	ng/ul	99
16) Anthracene	17.22	178	28747	0.537	ng/ul	100
19) Fluoranthene	19.15	202	20749	0.336	ng/ul	99
20) Pyrene	19.51	202	26613	0.423	ng/ul	99
21) Benzo(a)anthracene	21.27	228	29294	0.514	ng/ul	99
22) Chrysene	21.32	228	32216	0.544	ng/ul	100
24) Benzo(b)fluoranthene	22.85	252	49540	0.844	ng/ul	98
25) Benzo(k)fluoranthene	22.89	252	23609	0.396	ng/ul	99
26) Benzo(a)pyrene	23.42	252	18883	0.388	ng/ul	97
28) Dibenzo(a,h)anthracene	25.78	278	49198	0.920	ng/ul	98
29) Benzo(a,h,i)perylene	26.45	276	21825	0.387	ng/ul	99

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

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