

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080422\
 Data File : BN021009.D
 Acq On : 05 Aug 2022 03:32
 Operator : CG/JU
 Sample : SP5663
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SP5663

Quant Time: Aug 05 04:21:49 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 03 16:31:11 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.809	152	2480	0.400	ng/ul	0.00
4) Naphthalene-d8	10.596	136	7928	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.457	164	4462	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.205	188	8685	0.400	ng/ul	0.00
17) Chrysene-d12	21.409	240	7376	0.400	ng/ul	0.00
23) Perylene-d12	23.754	264	5988	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	2460	0.775	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.202	152	5014	0.432	ng/ul	0.00
18) Fluoranthene-d10	19.245	212	9757	0.428	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.240	88	12091	3.699	ng/ul#	80
5) Naphthalene	10.646	128	16381	0.673	ng/ul	98
7) 2-Methylnaphthalene	12.273	142	10138	0.686	ng/ul	99
8) 1-Methylnaphthalene	12.493	142	10887	0.737	ng/ul	99
10) Acenaphthylene	14.174	152	14190	0.680	ng/ul	98
11) Acenaphthene	14.517	153	11429	0.662	ng/ul	99
12) Fluorene	15.507	166	12613	0.669	ng/ul	98
14) Pentachlorophenol	16.872	266	1929	1.440	ng/ul	99
15) Phenanthrene	17.248	178	20172	0.649	ng/ul	99
16) Anthracene	17.340	178	17028	0.708	ng/ul	99
19) Fluoranthene	19.277	202	21864	0.670	ng/ul	98
20) Pyrene	19.640	202	22532	0.669	ng/ul	99
21) Benzo(a)anthracene	21.395	228	17288	0.718	ng/ul	100
22) Chrysene	21.447	228	20713	0.654	ng/ul	99
24) Benzo(b)fluoranthene	23.040	252	19006	0.686	ng/ul	96
25) Benzo(k)fluoranthene	23.087	252	19700	0.673	ng/ul	96
26) Benzo(a)pyrene	23.648	252	17212	0.681	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.169	276	18960	0.628	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.182	278	15249	0.639	ng/ul	95
29) Benzo(g,h,i)perylene	26.907	276	16653	0.618	ng/ul	99

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

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