

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN052422\
 Data File : BN019957.D
 Acq On : 25 May 2022 18:33
 Operator : CG/JU
 Sample : N2883-05MS
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EW4K6MS

Quant Time: May 26 01:15:53 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN052322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 23 14:58:28 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	4251	0.400	ng/ul	0.00
4) Naphthalene-d8	10.639	136	14569	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.474	164	9391	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.213	188	17401	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.403	240	12187	0.400	ng/ul	# 0.00
23) Perylene-d12	23.741	264	10317	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1825	0.460	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	6940	0.292	ng/ul	0.00
18) Fluoranthene-d10	19.239	212	15896	0.365	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.353	88	4920	1.189	ng/ul#	80
5) Naphthalene	10.689	128	12021	0.263	ng/ul	99
7) 2-Methylnaphthalene	12.300	142	10169	0.336	ng/ul#	100
8) 1-Methylnaphthalene	12.520	142	9712	0.345	ng/ul#	100
10) Acenaphthylene	14.192	152	12901	0.241	ng/ul	93
11) Acenaphthene	14.534	153	9196	0.254	ng/ul	96
12) Fluorene	15.520	166	13104	0.323	ng/ul#	87
15) Phenanthrene	17.255	178	41573	0.679	ng/ul#	91
16) Anthracene	17.344	178	16353	0.285	ng/ul#	81
19) Fluoranthene	19.272	202	26206	0.431	ng/ul#	93
20) Pyrene	19.634	202	33503	0.543	ng/ul#	95
21) Benzo(a)anthracene	21.385	228	16834	0.338	ng/ul#	69
22) Chrysene	21.438	228	24549	0.488	ng/ul#	74
24) Benzo(b)fluoranthene	23.031	252	16004	0.351	ng/ul#	1
25) Benzo(k)fluoranthene	23.080	252	12918	0.291	ng/ul#	1
26) Benzo(a)pyrene	23.639	252	14275	0.331	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.144	276	13847	0.306	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.165	278	10875	0.293	ng/ul#	54
29) Benzo(g,h,i)perylene	26.882	276	9385	0.243	ng/ul#	68

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

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