

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072522\
 Data File : BN020839.D
 Acq On : 26 Jul 2022 00:21
 Operator : CG/JU
 Sample : N3801-09MSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GBHX2MSD

Quant Time: Jul 26 00:50:06 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN072522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 25 18:06:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.825	152	2070	0.400	ng/ul	0.00
4) Naphthalene-d8	10.617	136	7465	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.470	164	4369	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.221	188	9319	0.400	ng/ul	0.00
17) Chrysene-d12	21.420	240	8462	0.400	ng/ul	-0.01
23) Perylene-d12	23.767	264	6968	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	642	0.258	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.218	152	4180	0.369	ng/ul	0.00
18) Fluoranthene-d10	19.253	212	11932	0.441	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.252	88	2505	0.980	ng/ul#	86
5) Naphthalene	10.667	128	7820	0.334	ng/ul	99
7) 2-Methylnaphthalene	12.289	142	4777	0.323	ng/ul	99
8) 1-Methylnaphthalene	12.509	142	5145	0.359	ng/ul	99
10) Acenaphthylene	14.187	152	7074	0.326	ng/ul	99
11) Acenaphthene	14.530	153	5640	0.321	ng/ul	99
12) Fluorene	15.524	166	6521	0.338	ng/ul	100
14) Pentachlorophenol	16.884	266	1227	0.874	ng/ul	97
15) Phenanthrene	17.264	178	11808	0.346	ng/ul	99
16) Anthracene	17.356	178	9568	0.350	ng/ul	99
19) Fluoranthene	19.285	202	14778	0.373	ng/ul	99
20) Pyrene	19.648	202	15251	0.369	ng/ul	99
21) Benzo(a)anthracene	21.406	228	10833	0.387	ng/ul	100
22) Chrysene	21.458	228	13497	0.375	ng/ul	99
24) Benzo(b)fluoranthene	23.054	252	12493	0.378	ng/ul	97
25) Benzo(k)fluoranthene	23.101	252	12441	0.364	ng/ul	99
26) Benzo(a)pyrene	23.665	252	10709	0.352	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.201	276	11800	0.368	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.218	278	9244	0.379	ng/ul	99
29) Benzo(g,h,i)perylene	26.933	276	10807	0.374	ng/ul	99

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

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