

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071923\
 Data File : BN026595.D
 Acq On : 19 Jul 2023 12:47
 Operator : MA/JU
 Sample : 03606-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR-04-293-308-071223

Quant Time: Jul 19 15:32:28 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 19 15:31:39 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	2799	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.737	136	8284	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.571	164	5529	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.322	188	12965	0.400	ng/ul	0.00
17) Chrysene-d12	21.524	240	12296	0.400	ng/ul	# 0.00
23) Perylene-d12	23.962	264	13275	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/ul	
6) 2-Methylnaphthalene-d10	12.327	152	3074	0.273	ng/ul	-0.01
18) Fluoranthene-d10	19.355	212	11472	0.269	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.318	88	10472	3.002	ng/ul#	75
5) Naphthalene	10.787	128	468	0.021	ng/ul#	51
7) 2-Methylnaphthalene	12.404	142	132	0.010	ng/ul#	59
8) 1-Methylnaphthalene	12.623	142	130	0.009	ng/ul	82
10) Acenaphthylene	14.294	152	149	0.006	ng/ul#	1
11) Acenaphthene	14.650	153	211	0.010	ng/ul#	72
12) Fluorene	15.626	166	104	0.005	ng/ul#	60
14) Pentachlorophenol	16.989	266	92	0.028	ng/ul#	84
15) Phenanthrene	17.365	178	891	0.022	ng/ul#	72
16) Anthracene	17.462	178	225	0.006	ng/ul#	10
19) Fluoranthene	19.383	202	926	0.016	ng/ul#	33
20) Pyrene	19.750	202	732	0.012	ng/ul#	37
21) Benzo(a)anthracene	21.509	228	466	0.010	ng/ul#	31
22) Chrysene	21.559	228	412	0.008	ng/ul#	39
24) Benzo(b)fluoranthene	23.228	252	430	0.008	ng/ul#	1
25) Benzo(k)fluoranthene	23.272	252	210	0.004	ng/ul#	1
26) Benzo(a)pyrene	23.857	252	24	0.001	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.525	276	258	0.004	ng/ul#	54
28) Dibenzo(a,h)anthracene	26.522	278	185	0.004	ng/ul#	1
29) Benzo(g,h,i)perylene	27.283	276	161	0.003	ng/ul#	1

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

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