

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071923\
 Data File : BN026596.D
 Acq On : 19 Jul 2023 13:24
 Operator : MA/JU
 Sample : 03606-03
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

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

Quant Time: Jul 19 14:35:12 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 09:49:57 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	2615	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.737	136	8014	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.571	164	5384	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.323	188	12367	0.400	ng/ul	-0.01
17) Chrysene-d12	21.524	240	11566	0.400	ng/ul	# 0.00
23) Perylene-d12	23.962	264	12414	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	3769	0.346	ng/ul	-0.02
18) Fluoranthene-d10	19.355	212	12462	0.311	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.318	88	12703	3.898	ng/ul#	82
5) Naphthalene	10.787	128	623	0.028	ng/ul#	60
7) 2-Methylnaphthalene	12.404	142	189	0.015	ng/ul	89
8) 1-Methylnaphthalene	12.624	142	163	0.012	ng/ul#	79
10) Acenaphthylene	14.298	152	165	0.006	ng/ul#	1
11) Acenaphthene	14.650	153	236	0.012	ng/ul#	70
12) Fluorene	15.626	166	120	0.006	ng/ul#	64
14) Pentachlorophenol	16.993	266	23	0.007	ng/ul#	67
15) Phenanthrene	17.365	178	970	0.025	ng/ul#	73
16) Anthracene	17.458	178	192	0.006	ng/ul#	15
19) Fluoranthene	19.388	202	1099	0.020	ng/ul#	41
20) Pyrene	19.750	202	720	0.012	ng/ul#	39
21) Benzo(a)anthracene	21.509	228	440	0.010	ng/ul#	31
22) Chrysene	21.565	228	482	0.010	ng/ul#	42
24) Benzo(b)fluoranthene	23.222	252	415	0.009	ng/ul#	1
25) Benzo(k)fluoranthene	23.272	252	190	0.004	ng/ul#	1
26) Benzo(a)pyrene	23.851	252	121	0.003	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.518	276	260	0.005	ng/ul#	49
28) Dibenzo(a,h)anthracene	26.522	278	169	0.004	ng/ul#	1
29) Benzo(g,h,i)perylene	27.293	276	210	0.004	ng/ul#	1

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

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