

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071923\
 Data File : BN026597.D
 Acq On : 19 Jul 2023 15:20
 Operator : MA/JU
 Sample : 03635-02
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 GWBR-04-359-374-071423

Quant Time: Jul 19 15:55:00 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 15:54:02 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	1433	0.400	ng/ul	-0.07
4) Naphthalene-d8	10.677	136	4751	0.400	ng/ul	-0.07
9) Acenaphthene-d10	14.515	164	3017	0.400	ng/ul	-0.06
13) Phenanthrene-d10	17.267	188	7815	0.400	ng/ul	#-0.06
17) Chrysene-d12	21.479	240	12725	0.400	ng/ul	#-0.05
23) Perylene-d12	23.905	264	15317	0.400	ng/ul	#-0.06
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/ul	
6) 2-Methylnaphthalene-d10	12.272	152	816	0.126	ng/ul	-0.07
18) Fluoranthene-d10	19.294	212	5268	0.119	ng/ul	-0.06
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.297	88	595	0.333	ng/ul#	1
5) Naphthalene	10.732	128	30	0.002	ng/ul#	1
7) 2-Methylnaphthalene	12.343	142	25	0.003	ng/ul	81
8) 1-Methylnaphthalene	12.530	142	48	0.006	ng/ul#	31
10) Acenaphthylene	14.237	152	196	0.013	ng/ul#	1
11) Acenaphthene	14.584	153	32	0.003	ng/ul#	54
12) Fluorene	15.570	166	39	0.003	ng/ul#	48
14) Pentachlorophenol	16.954	266	13	0.007	ng/ul#	46
15) Phenanthrene	17.309	178	122	0.005	ng/ul#	1
16) Anthracene	17.398	178	123	0.006	ng/ul#	1
19) Fluoranthene	19.326	202	395	0.007	ng/ul#	21
20) Pyrene	19.693	202	141	0.002	ng/ul#	1
21) Benzo(a)anthracene	21.471	228	228	0.005	ng/ul#	12
22) Chrysene	21.520	228	200	0.004	ng/ul#	16
24) Benzo(b)fluoranthene	23.180	252	208	0.003	ng/ul#	1
25) Benzo(k)fluoranthene	23.218	252	18	0.000	ng/ul#	1
26) Benzo(a)pyrene	23.797	252	94	0.002	ng/ul#	1
27) Indeno(1,2,3-cd)pyrene	26.427	276	29	0.000	ng/ul#	67
28) Dibenzo(a,h)anthracene	26.447	278	37	0.001	ng/ul#	1
29) Benzo(g,h,i)perylene	27.235	276	8	0.000	ng/ul#	1

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

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