

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026431.D
 Acq On : 09 Jul 2023 20:20
 Operator : MA/JU
 Sample : 03351-02MS
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW10MS

Quant Time: Jul 10 01:23:44 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	5827	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	24552	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	13167	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.322	188	19114	0.400	ng/ul	0.00
17) Chrysene-d12	21.511	240	11641	0.400	ng/ul	# 0.00
23) Perylene-d12	23.967	264	13363	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.306	96	975	0.149	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.326	152	6333	0.168	ng/ul	0.00
18) Fluoranthene-d10	19.349	212	9013	0.203	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.339	88	2537	0.364	ng/ul	96
5) Naphthalene	10.781	128	13358	0.190	ng/ul	99
7) 2-Methylnaphthalene	12.398	142	7194	0.161	ng/ul	99
8) 1-Methylnaphthalene	12.612	142	7889	0.170	ng/ul	99
10) Acenaphthylene	14.288	152	11389	0.185	ng/ul	99
11) Acenaphthene	14.626	153	10258	0.203	ng/ul	95
12) Fluorene	15.620	166	9195	0.159	ng/ul	99
14) Pentachlorophenol	16.984	266	1564	0.228	ng/ul	99
15) Phenanthrene	17.360	178	12861	0.204	ng/ul	99
16) Anthracene	17.457	178	10737	0.191	ng/ul	95
19) Fluoranthene	19.377	202	11705	0.188	ng/ul	97
20) Pyrene	19.740	202	12499	0.193	ng/ul#	92
21) Benzo(a)anthracene	21.497	228	10160	0.212	ng/ul	100
22) Chrysene	21.546	228	11031	0.221	ng/ul	99
24) Benzo(b)fluoranthene	23.215	252	11483	0.201	ng/ul	93
25) Benzo(k)fluoranthene	23.259	252	12223	0.213	ng/ul	94
26) Benzo(a)pyrene	23.864	252	10673	0.215	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.537	276	13421	0.248	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.551	278	10360	0.253	ng/ul	93
29) Benzo(g,h,i)perylene	27.332	276	11919	0.248	ng/ul	97

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

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