

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072023\
 Data File : BN026636.D
 Acq On : 21 Jul 2023 20:09
 Operator : MA/JU
 Sample : 03536-02MS
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW57MS

Quant Time: Jul 22 00:59:04 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 : Fri Jul 21 19:02:16 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	2741	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.630	136	9171	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.472	164	5495	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.219	188	11688	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.411	240	8658	0.400	ng/ul	#-0.01
23) Perylene-d12	23.778	264	8639	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.204	96	460	0.127	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.225	152	1522	0.122	ng/ul	-0.02
18) Fluoranthene-d10	19.250	212	3912	0.130	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.237	88	2011	0.589	ng/ul#	1
5) Naphthalene	10.680	128	4652	0.185	ng/ul#	93
7) 2-Methylnaphthalene	12.297	142	2397	0.168	ng/ul	99
8) 1-Methylnaphthalene	12.517	142	2242	0.146	ng/ul	92
10) Acenaphthylene	14.194	152	2943	0.109	ng/ul#	89
11) Acenaphthene	14.532	153	2646	0.129	ng/ul	99
12) Fluorene	15.526	166	2579	0.120	ng/ul#	96
14) Pentachlorophenol	16.886	266	434	0.147	ng/ul	98
15) Phenanthrene	17.261	178	5820	0.156	ng/ul	96
16) Anthracene	17.359	178	3770	0.117	ng/ul#	90
19) Fluoranthene	19.278	202	8513	0.210	ng/ul#	90
20) Pyrene	19.641	202	10630	0.241	ng/ul#	90
21) Benzo(a)anthracene	21.393	228	6254	0.197	ng/ul#	73
22) Chrysene	21.446	228	6817	0.188	ng/ul#	76
24) Benzo(b)fluoranthene	23.056	252	8655	0.257	ng/ul#	24
25) Benzo(k)fluoranthene	23.103	252	5281	0.150	ng/ul#	24
26) Benzo(a)pyrene	23.673	252	4746	0.158	ng/ul#	10
27) Indeno(1,2,3-cd)pyrene	26.231	276	6161	0.157	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.241	278	4309	0.143	ng/ul#	30
29) Benzo(g,h,i)perylene	26.992	276	919	0.027	ng/ul#	1

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

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