

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026433.D
 Acq On : 09 Jul 2023 21:33
 Operator : MA/JU
 Sample : 03349-07MS
 Misc :
 ALS Vial : 61 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW01MS

Quant Time: Jul 10 01:37:07 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	7.921	152	5226	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.730	136	20405	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.565	164	11867	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.321	188	17819	0.400	ng/ul	0.00
17) Chrysene-d12	21.513	240	11343	0.400	ng/ul	0.00
23) Perylene-d12	23.954	264	11829	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	791	0.134	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.325	152	5209	0.166	ng/ul	0.00
18) Fluoranthene-d10	19.353	212	9854	0.228	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	1944	0.311	ng/ul	97
5) Naphthalene	10.780	128	11033	0.189	ng/ul	98
7) 2-Methylnaphthalene	12.402	142	5937	0.159	ng/ul	100
8) 1-Methylnaphthalene	12.611	142	6670	0.173	ng/ul	100
10) Acenaphthylene	14.287	152	10178	0.183	ng/ul	99
11) Acenaphthene	14.630	153	9112	0.200	ng/ul	94
12) Fluorene	15.620	166	8469	0.163	ng/ul	99
14) Pentachlorophenol	16.987	266	1350	0.211	ng/ul	99
15) Phenanthrene	17.363	178	12067	0.205	ng/ul	98
16) Anthracene	17.460	178	9977	0.191	ng/ul	95
19) Fluoranthene	19.381	202	13716	0.226	ng/ul	100
20) Pyrene	19.744	202	14617	0.232	ng/ul	97
21) Benzo(a)anthracene	21.499	228	9971	0.213	ng/ul	99
22) Chrysene	21.551	228	11041	0.227	ng/ul	98
24) Benzo(b)fluoranthene	23.211	252	10853	0.215	ng/ul	89
25) Benzo(k)fluoranthene	23.261	252	11747	0.232	ng/ul#	91
26) Benzo(a)pyrene	23.852	252	9815	0.224	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.513	276	12396	0.259	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.526	278	9448	0.261	ng/ul	93
29) Benzo(g,h,i)perylene	27.294	276	11198	0.263	ng/ul	98

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

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