

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070823\
 Data File : BN026434.D
 Acq On : 09 Jul 2023 22:10
 Operator : MA/JU
 Sample : 03349-08MSD
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW01MSD

Quant Time: Jul 10 01:38:37 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	7.930	152	4335	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	15100	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.566	164	8483	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.322	188	12128	0.400	ng/ul	0.00
17) Chrysene-d12	21.511	240	8066	0.400	ng/ul	0.00
23) Perylene-d12	23.958	264	8473	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	903	0.185	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.327	152	5577	0.240	ng/ul	0.00
18) Fluoranthene-d10	19.350	212	10149	0.331	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.335	88	2378	0.459	ng/ul	100
5) Naphthalene	10.781	128	11916	0.275	ng/ul	99
7) 2-Methylnaphthalene	12.398	142	6291	0.228	ng/ul	99
8) 1-Methylnaphthalene	12.613	142	7142	0.250	ng/ul	99
10) Acenaphthylene	14.288	152	10563	0.266	ng/ul	98
11) Acenaphthene	14.626	153	9539	0.294	ng/ul	95
12) Fluorene	15.620	166	8660	0.232	ng/ul	100
14) Pentachlorophenol	16.988	266	1454	0.334	ng/ul	100
15) Phenanthrene	17.364	178	12078	0.301	ng/ul	98
16) Anthracene	17.461	178	9914	0.278	ng/ul	96
19) Fluoranthene	19.382	202	14058	0.326	ng/ul	98
20) Pyrene	19.745	202	15005	0.335	ng/ul	97
21) Benzo(a)anthracene	21.494	228	10382	0.312	ng/ul	98
22) Chrysene	21.547	228	11904	0.344	ng/ul	99
24) Benzo(b)fluoranthene	23.210	252	11395	0.315	ng/ul	90
25) Benzo(k)fluoranthene	23.259	252	12499	0.344	ng/ul#	92
26) Benzo(a)pyrene	23.850	252	9918	0.316	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	26.527	276	13106	0.382	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.538	278	9940	0.383	ng/ul	97
29) Benzo(g,h,i)perylene	27.316	276	11794	0.386	ng/ul	98

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

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