

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM052523\
 Data File : BM040031.D
 Acq On : 26 May 2023 09:17
 Operator : CG/JU
 Sample : 02861-08MSD
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 H0FL1MSD

Quant Time: May 26 11:31:13 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM052523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 25 15:01:16 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.012	152	7860	0.400	ng/ul	0.00
4) Naphthalene-d8	10.826	136	25530	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.645	164	12394	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.390	188	23528	0.400	ng/ul	0.00
17) Chrysene-d12	21.566	240	11752	0.400	ng/ul	0.00
23) Perylene-d12	24.013	264	10257	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.384	96	1918	0.210	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.410	152	9965	0.319	ng/ul	0.00
18) Fluoranthene-d10	19.411	212	17007	0.481	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.417	88	4903	0.509	ng/ul	93
5) Naphthalene	10.876	128	19823	0.318	ng/ul	100
7) 2-Methylnaphthalene	12.481	142	11873	0.315	ng/ul	100
8) 1-Methylnaphthalene	12.701	142	12255	0.313	ng/ul	99
10) Acenaphthylene	14.368	152	16286	0.330	ng/ul	99
11) Acenaphthene	14.710	153	13603	0.338	ng/ul	99
12) Fluorene	15.691	166	15043	0.359	ng/ul	99
14) Pentachlorophenol	17.035	266	3493	0.732	ng/ul	98
15) Phenanthrene	17.428	178	24974	0.375	ng/ul	100
16) Anthracene	17.521	178	17694	0.317	ng/ul	99
19) Fluoranthene	19.439	202	24188	0.457	ng/ul	99
20) Pyrene	19.801	202	23463	0.438	ng/ul	99
21) Benzo(a)anthracene	21.548	228	16033	0.405	ng/ul	99
22) Chrysene	21.604	228	17488	0.416	ng/ul	100
24) Benzo(b)fluoranthene	23.264	252	16356	0.402	ng/ul	98
25) Benzo(k)fluoranthene	23.314	252	16175	0.394	ng/ul	99
26) Benzo(a)pyrene	23.904	252	11249	0.330	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	26.556	276	16976	0.346	ng/ul	100
28) Dibenzo(a,h)anthracene	26.577	278	13794	0.348	ng/ul	100
29) Benzo(g,h,i)perylene	27.335	276	13774	0.343	ng/ul	99

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

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