

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060121\
 Data File : BM030259.D
 Acq On : 02 Jun 2021 14:56
 Operator : CG/JU
 Sample : M2528-03MSD
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YARK5MSD

Quant Time: Jun 02 15:26:47 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM060121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 02 02:39:12 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.864	152	4991	0.40	ng/ul	0.00
4) Naphthalene-d8	10.653	136	17541	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.485	164	9334	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.218	188	17862	0.40	ng/ul	0.00
17) Chrysene-d12	21.377	240	11332	0.40	ng/ul	0.00
23) Perylene-d12	23.663	264	9212	0.40	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	1357	0.28	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.236	152	8697	0.35	ng/ul	0.00
18) Fluoranthene-d10	19.237	212	16557	0.43	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.369	88	2148	0.41	ng/ul	96
5) Naphthalene	10.703	128	17731	0.34	ng/ul	99
7) 2-Methylnaphthalene	12.313	142	11860	0.34	ng/ul	99
8) 1-Methylnaphthalene	12.533	142	10971	0.33	ng/ul	99
10) Acenaphthylene	14.205	152	14058	0.32	ng/ul	100
11) Acenaphthene	14.546	153	12065	0.32	ng/ul	99
12) Fluorene	15.528	166	13481	0.31	ng/ul	99
14) Pentachlorophenol	16.867	266	3302	0.59	ng/ul	98
15) Phenanthrene	17.259	178	22350	0.35	ng/ul	100
16) Anthracene	17.350	178	18871	0.34	ng/ul	100
19) Fluoranthene	19.264	202	23923	0.41	ng/ul	100
20) Pyrene	19.626	202	24077	0.41	ng/ul	99
21) Benzo(a)anthracene	21.360	228	17696	0.39	ng/ul	99
22) Chrysene	21.410	228	19176	0.39	ng/ul	100
24) Benzo(b)fluoranthene	22.970	252	18534	0.40	ng/ul	99
25) Benzo(k)fluoranthene	23.016	252	17288	0.40	ng/ul	100
26) Benzo(a)pyrene	23.559	252	14310	0.39	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.982	276	18589	0.38	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.996	278	15135	0.38	ng/ul	99
29) Benzo(g,h,i)perylene	26.694	276	16290	0.38	ng/ul	99

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

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