

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM060121\
 Data File : BM030248.D
 Acq On : 01 Jun 2021 22:11
 Operator : CG/JU
 Sample : M2528-03MSD
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 YARK5MSD

Quant Time: Jun 02 01:11:26 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 01:10:46 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	4667	0.40	ng/ul	0.00
4) Naphthalene-d8	10.657	136	16195	0.40	ng/ul	0.00
9) Acenaphthene-d10	14.485	164	8927	0.40	ng/ul	0.00
13) Phenanthrene-d10	17.218	188	18248	0.40	ng/ul	0.00
17) Chrysene-d12	21.374	240	14466	0.40	ng/ul	0.00
23) Perylene-d12	23.661	264	12047	0.40	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.338	96	455	0.10	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.241	152	7746	0.33	ng/ul	0.00
18) Fluoranthene-d10	19.237	212	18729	0.39	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.369	88	1916	0.39	ng/ul#	66
5) Naphthalene	10.702	128	17409	0.36	ng/ul	99
7) 2-Methylnaphthalene	12.313	142	11741	0.36	ng/ul	100
8) 1-Methylnaphthalene	12.533	142	10873	0.35	ng/ul	100
10) Acenaphthylene	14.208	152	14364	0.34	ng/ul	99
11) Acenaphthene	14.550	153	12356	0.34	ng/ul	100
12) Fluorene	15.532	166	14247	0.35	ng/ul	100
14) Pentachlorophenol	16.867	266	3713	0.65	ng/ul	98
15) Phenanthrene	17.259	178	24973	0.38	ng/ul	99
16) Anthracene	17.349	178	21266	0.38	ng/ul	99
19) Fluoranthene	19.268	202	30230	0.41	ng/ul	99
20) Pyrene	19.626	202	30978	0.41	ng/ul	98
21) Benzo(a)anthracene	21.362	228	25331	0.44	ng/ul	100
22) Chrysene	21.413	228	26939	0.43	ng/ul	100
24) Benzo(b)fluoranthene	22.970	252	26203	0.44	ng/ul	97
25) Benzo(k)fluoranthene	23.016	252	25232	0.44	ng/ul	96
26) Benzo(a)pyrene	23.559	252	20399	0.42	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.982	276	24675	0.39	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.996	278	19984	0.39	ng/ul	99
29) Benzo(g,h,i)perylene	26.691	276	21161	0.38	ng/ul	99

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

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