

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022824\
 Data File : BN030153.D
 Acq On : 29 Feb 2024 05:51
 Operator : MA/JU
 Sample : P1494-16MSD
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BH3W6MSD

Quant Time: Feb 29 06:17:44 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 27 23:49:29 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.876	152	1969	0.400	ng/ul	0.02
4) Naphthalene-d8	10.672	136	5464	0.400	ng/ul	0.02
9) Acenaphthene-d10	14.517	164	2754	0.400	ng/ul	0.02
13) Phenanthrene-d10	17.269	188	5376	0.400	ng/ul	0.00
17) Chrysene-d12	21.471	240	5313	0.400	ng/ul	0.00
23) Perylene-d12	23.847	264	4735	0.400	ng/ul	# 0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	943	0.384	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.267	152	3046	0.438	ng/ul	0.02
18) Fluoranthene-d10	19.300	212	7670	0.439	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.361	88	2521	0.943	ng/ul	99
5) Naphthalene	10.722	128	6251	0.407	ng/ul	98
7) 2-Methylnaphthalene	12.339	142	3624	0.411	ng/ul	99
8) 1-Methylnaphthalene	12.558	142	3863	0.430	ng/ul	99
10) Acenaphthylene	14.239	152	5431	0.404	ng/ul	97
11) Acenaphthene	14.577	153	4168	0.408	ng/ul	97
12) Fluorene	15.571	166	4427	0.414	ng/ul	97
14) Pentachlorophenol	16.918	266	1288	1.074	ng/ul#	51
15) Phenanthrene	17.311	178	7407	0.441	ng/ul	95
16) Anthracene	17.408	178	6141	0.435	ng/ul#	95
19) Fluoranthene	19.333	202	9574	0.392	ng/ul#	94
20) Pyrene	19.695	202	10464	0.418	ng/ul#	92
21) Benzo(a)anthracene	21.456	228	7463	0.390	ng/ul	98
22) Chrysene	21.506	228	8638	0.412	ng/ul	99
24) Benzo(b)fluoranthene	23.125	252	7948	0.446	ng/ul	97
25) Benzo(k)fluoranthene	23.172	252	9206	0.466	ng/ul	95
26) Benzo(a)pyrene	23.745	252	7281	0.435	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.320	276	8281	0.407	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.363	278	6327	0.397	ng/ul#	86
29) Benzo(g,h,i)perylene	27.057	276	7601	0.412	ng/ul	93

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

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