

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030228.D
 Acq On : 03 Mar 2024 21:11
 Operator : MA/JU
 Sample : P1501-04MSD
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E10A0MSD

Quant Time: Mar 03 22:55:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN030224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 02 16:09:30 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	1400	0.400	ng/ul	0.00
4) Naphthalene-d8	10.634	136	3690	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.479	164	1730	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.239	188	3251	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.445	240	2971	0.400	ng/ul	# 0.00
23) Perylene-d12	23.801	264	3434	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.311	96	462	0.238	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.228	152	1795	0.376	ng/ul	-0.01
18) Fluoranthene-d10	19.273	212	3840	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.340	88	1419	0.675	ng/ul	93
5) Naphthalene	10.683	128	3883	0.357	ng/ul	98
7) 2-Methylnaphthalene	12.305	142	2149	0.364	ng/ul	99
8) 1-Methylnaphthalene	12.520	142	2284	0.365	ng/ul	99
10) Acenaphthylene	14.202	152	3467	0.358	ng/ul	98
11) Acenaphthene	14.540	153	2496	0.352	ng/ul	97
12) Fluorene	15.539	166	2674	0.370	ng/ul	95
14) Pentachlorophenol	16.889	266	921	0.993	ng/ul	98
15) Phenanthrene	17.282	178	4045	0.371	ng/ul	99
16) Anthracene	17.374	178	3678	0.384	ng/ul	99
19) Fluoranthene	19.301	202	5045	0.382	ng/ul	97
20) Pyrene	19.668	202	5353	0.373	ng/ul	95
21) Benzo(a)anthracene	21.430	228	4446	0.423	ng/ul	99
22) Chrysene	21.480	228	5133	0.407	ng/ul	99
24) Benzo(b)fluoranthene	23.084	252	5369	0.405	ng/ul	95
25) Benzo(k)fluoranthene	23.131	252	5949	0.381	ng/ul#	93
26) Benzo(a)pyrene	23.693	252	5184	0.385	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.226	276	7077	0.411	ng/ul	95
28) Dibenzo(a,h)anthracene	26.260	278	5587	0.419	ng/ul	93
29) Benzo(g,h,i)perylene	26.967	276	6318	0.396	ng/ul#	93

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

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