

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030216.D
 Acq On : 03 Mar 2024 13:22
 Operator : MA/JU
 Sample : P1512-05MSD
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 E0S96MSD

Quant Time: Mar 03 22:52:53 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	1776	0.400	ng/ul	0.00
4) Naphthalene-d8	10.628	136	4554	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.479	164	2133	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.235	188	4040	0.400	ng/ul	-0.01
17) Chrysene-d12	21.444	240	3617	0.400	ng/ul	# 0.00
23) Perylene-d12	23.797	264	4385	0.400	ng/ul	#-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1269	0.516	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.229	152	1712	0.291	ng/ul	-0.01
18) Fluoranthene-d10	19.272	212	3376	0.286	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.340	88	3576	1.340	ng/ul#	90
5) Naphthalene	10.678	128	5783	0.430	ng/ul	98
7) 2-Methylnaphthalene	12.300	142	2224	0.305	ng/ul	100
8) 1-Methylnaphthalene	12.515	142	2281	0.295	ng/ul	100
10) Acenaphthylene	14.201	152	3280	0.275	ng/ul	98
11) Acenaphthene	14.539	153	2421	0.277	ng/ul	98
12) Fluorene	15.538	166	2573	0.289	ng/ul	99
14) Pentachlorophenol	16.884	266	467	0.405	ng/ul	97
15) Phenanthrene	17.277	178	4375	0.323	ng/ul	99
16) Anthracene	17.370	178	3430	0.288	ng/ul	100
19) Fluoranthene	19.300	202	5332	0.332	ng/ul	99
20) Pyrene	19.662	202	5486	0.314	ng/ul	99
21) Benzo(a)anthracene	21.427	228	4331	0.339	ng/ul	99
22) Chrysene	21.479	228	4640	0.302	ng/ul	97
24) Benzo(b)fluoranthene	23.084	252	5270	0.312	ng/ul	96
25) Benzo(k)fluoranthene	23.131	252	5138	0.258	ng/ul	96
26) Benzo(a)pyrene	23.695	252	5064	0.294	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.232	276	6717	0.306	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.256	278	5134	0.302	ng/ul	96
29) Benzo(g,h,i)perylene	26.970	276	6034	0.296	ng/ul	96

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

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