

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021323\
 Data File : BN024011.D
 Acq On : 13 Feb 2023 13:16
 Operator : CG/JU
 Sample : 01443-08
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3E97

Quant Time: Feb 14 01:22:44 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 14 01:13:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	3701	0.400	ng/u1	0.00
4) Naphthalene-d8	10.639	136	9656	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.488	164	5617	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.238	188	12598	0.400	ng/u1	# 0.00
17) Chrysene-d12	21.432	240	11880	0.400	ng/u1	# 0.00
23) Perylene-d12	23.820	264	9255	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.231	96	1578	0.398	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.234	152	3771	0.250	ng/u1	0.00
18) Fluoranthene-d10	19.272	212	11223	0.352	ng/u1	0.00
Target Compounds						
						Qvalue
5) Naphthalene	10.689	128	7284	0.276	ng/u1	100
7) 2-Methylnaphthalene	12.306	142	4793	0.279	ng/u1	100
10) Acenaphthylene	14.210	152	12381	0.576	ng/u1	98
12) Fluorene	15.538	166	6855	0.300	ng/u1	97
14) Pentachlorophenol	16.892	266	1507	0.363	ng/u1	99
15) Phenanthrene	17.280	178	21554	0.574	ng/u1	99
16) Anthracene	17.373	178	7119	0.226	ng/u1	98
19) Fluoranthene	19.299	202	18930	0.479	ng/u1#	94
20) Pyrene	19.667	202	22372	0.547	ng/u1	98
21) Benzo(a)anthracene	21.417	228	30694	0.761	ng/u1	100
22) Chrysene	21.470	228	18096	0.426	ng/u1	99
24) Benzo(b)fluoranthene	23.092	252	13880	0.339	ng/u1	92
25) Benzo(k)fluoranthene	23.145	252	21586	0.540	ng/u1#	98
27) Indeno(1,2,3-cd)pyrene	26.275	276	18637	0.390	ng/u1#	55
28) Dibenzo(a,h)anthracene	26.299	278	16441	0.430	ng/u1	98
29) Benzo(g,h,i)perylene	27.026	276	24774	0.630	ng/u1	97

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

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