

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022122\
 Data File : BN018664.D
 Acq On : 21 Feb 2022 13:45
 Operator : CG/JU
 Sample : N1486-03
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3943

Quant Time: Feb 21 14:13:02 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN021822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 18 13:39:55 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.913	152	6526	0.400	ng/ul	0.00
4) Naphthalene-d8	10.721	136	15831	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.552	164	7324	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.288	188	12898	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.468	240	10057	0.400	ng/ul	# 0.00
23) Perylene-d12	23.867	264	9251	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.284	96	4097	0.553	ng/ul	-0.01
6) 2-Methylnaphthalene-d10	12.310	152	6022	0.278	ng/ul	0.00
18) Fluoranthene-d10	19.317	212	9987	0.332	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.770	128	12768	0.285	ng/ul	99
7) 2-Methylnaphthalene	12.381	142	14030	0.485	ng/ul	100
10) Acenaphthylene	14.269	152	7288	0.221	ng/ul	98
12) Fluorene	15.593	166	12585	0.443	ng/ul	98
14) Pentachlorophenol	16.942	266	1592	0.605	ng/ul	99
15) Phenanthrene	17.330	178	12421	0.293	ng/ul	99
16) Anthracene	17.423	178	17228	0.455	ng/ul	99
19) Fluoranthene	19.345	202	11840	0.280	ng/ul#	94
20) Pyrene	19.707	202	15349	0.352	ng/ul#	94
21) Benzo(a)anthracene	21.453	228	19187	0.492	ng/ul	99
22) Chrysene	21.506	228	22104	0.545	ng/ul	100
24) Benzo(b)fluoranthene	23.136	252	35057	0.852	ng/ul	96
25) Benzo(k)fluoranthene	23.183	252	16351	0.398	ng/ul#	95
26) Benzo(a)pyrene	23.762	252	14717	0.405	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.370	278	33319	0.918	ng/ul#	95
29) Benzo(g,h,i)perylene	27.114	276	15599	0.394	ng/ul#	92

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

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