

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103122\
 Data File : BN022424.D
 Acq On : 01 Nov 2022 05:59
 Operator : CG/JU
 Sample : PB148647BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS647

Quant Time: Nov 01 06:30:34 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 01 01:09:32 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.928	152	3868	0.400	ng/ul	0.00
4) Naphthalene-d8	10.750	136	12649	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.600	164	7001	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.353	188	13074	0.400	ng/ul	0.00
17) Chrysene-d12	21.553	240	7400	0.400	ng/ul	0.00
23) Perylene-d12	23.996	264	5483	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	3175	0.632	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.351	152	6412	0.335	ng/ul	0.00
18) Fluoranthene-d10	19.389	212	9873	0.397	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.274	88	7956	1.541	ng/ul	88
5) Naphthalene	10.805	128	12180	0.326	ng/ul	100
7) 2-Methylnaphthalene	12.422	142	7566	0.321	ng/ul	100
8) 1-Methylnaphthalene	12.642	142	8182	0.338	ng/ul	100
10) Acenaphthylene	14.323	152	9828	0.313	ng/ul	99
11) Acenaphthene	14.660	153	8464	0.319	ng/ul	100
12) Fluorene	15.650	166	9153	0.297	ng/ul	99
14) Pentachlorophenol	17.007	266	941	0.303	ng/ul	99
15) Phenanthrene	17.395	178	13970	0.318	ng/ul	99
16) Anthracene	17.488	178	11374	0.308	ng/ul	98
19) Fluoranthene	19.417	202	12973	0.383	ng/ul	98
20) Pyrene	19.784	202	12890	0.379	ng/ul	99
21) Benzo(a)anthracene	21.535	228	9209	0.332	ng/ul	98
22) Chrysene	21.588	228	9974	0.339	ng/ul	99
24) Benzo(b)fluoranthene	23.245	252	9163	0.333	ng/ul	91
25) Benzo(k)fluoranthene	23.295	252	8702	0.325	ng/ul#	92
26) Benzo(a)pyrene	23.888	252	7650	0.349	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.561	276	10066	0.369	ng/ul#	40
28) Dibenzo(a,h)anthracene	26.581	278	7919	0.363	ng/ul	94
29) Benzo(g,h,i)perylene	27.356	276	8510	0.383	ng/ul	98

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

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