

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN103122\
 Data File : BN022429.D
 Acq On : 01 Nov 2022 09:42
 Operator : CG/JU
 Sample : PB148653BS
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS653

Quant Time: Nov 02 00:52:18 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	7.927	152	4062	0.400	ng/ul	0.00
4) Naphthalene-d8	10.754	136	13412	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.599	164	7471	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.356	188	14270	0.400	ng/ul	0.00
17) Chrysene-d12	21.553	240	8279	0.400	ng/ul	0.00
23) Perylene-d12	24.000	264	5670	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	3423	0.648	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.349	152	6997	0.345	ng/ul	0.00
18) Fluoranthene-d10	19.387	212	11219	0.404	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.274	88	8925	1.647	ng/ul#	87
5) Naphthalene	10.804	128	13508	0.341	ng/ul	100
7) 2-Methylnaphthalene	12.421	142	8346	0.334	ng/ul	100
8) 1-Methylnaphthalene	12.641	142	9032	0.352	ng/ul	100
10) Acenaphthylene	14.321	152	10743	0.320	ng/ul	100
11) Acenaphthene	14.659	153	9365	0.331	ng/ul	100
12) Fluorene	15.649	166	10153	0.309	ng/ul	98
14) Pentachlorophenol	17.010	266	954	0.281	ng/ul	99
15) Phenanthrene	17.394	178	15788	0.330	ng/ul	99
16) Anthracene	17.491	178	12814	0.317	ng/ul	99
19) Fluoranthene	19.420	202	15180	0.400	ng/ul	99
20) Pyrene	19.782	202	15038	0.395	ng/ul	98
21) Benzo(a)anthracene	21.535	228	10671	0.344	ng/ul	98
22) Chrysene	21.591	228	11534	0.351	ng/ul	99
24) Benzo(b)fluoranthene	23.245	252	9938	0.349	ng/ul	94
25) Benzo(k)fluoranthene	23.295	252	9660	0.349	ng/ul#	94
26) Benzo(a)pyrene	23.891	252	8226	0.362	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.565	276	10471	0.371	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.585	278	8292	0.368	ng/ul	96
29) Benzo(g,h,i)perylene	27.356	276	8884	0.386	ng/ul	98

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

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