

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030923\
 Data File : BN024192.D
 Acq On : 09 Mar 2023 06:17
 Operator : CG/JU
 Sample : PB151200BS
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS200

Quant Time: Mar 09 06:50:26 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 09 03:48:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	7628	0.400	ng/u1	0.00
4) Naphthalene-d8	10.872	136	21657	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.689	164	11185	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.430	188	22837	0.400	ng/u1	0.00
17) Chrysene-d12	21.621	240	16296	0.400	ng/u1	0.00
23) Perylene-d12	24.138	264	12565	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.406	96	5875	0.700	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.451	152	9828	0.382	ng/u1	0.00
18) Fluoranthene-d10	19.455	212	17432	0.402	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.443	88	14863	1.628	ng/u1	90
5) Naphthalene	10.922	128	20141	0.373	ng/u1	100
7) 2-Methylnaphthalene	12.528	142	12070	0.368	ng/u1	99
8) 1-Methylnaphthalene	12.742	142	12730	0.369	ng/u1	100
10) Acenaphthylene	14.407	152	16676	0.357	ng/u1	100
11) Acenaphthene	14.749	153	13234	0.368	ng/u1	99
12) Fluorene	15.735	166	14384	0.358	ng/u1	99
14) Pentachlorophenol	17.076	266	3713	0.585	ng/u1#	73
15) Phenanthrene	17.472	178	24056	0.368	ng/u1	100
16) Anthracene	17.565	178	19896	0.347	ng/u1	99
19) Fluoranthene	19.483	202	24893	0.384	ng/u1	97
20) Pyrene	19.845	202	24926	0.379	ng/u1	97
21) Benzo(a)anthracene	21.603	228	18452	0.347	ng/u1	99
22) Chrysene	21.659	228	21078	0.375	ng/u1	99
24) Benzo(b)fluoranthene	23.363	252	18629	0.388	ng/u1	98
25) Benzo(k)fluoranthene	23.419	252	18674	0.375	ng/u1	97
26) Benzo(a)pyrene	24.021	252	14795	0.358	ng/u1	96
27) Indeno(1,2,3-cd)pyrene	26.777	276	15822	0.340	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.804	278	11531	0.326	ng/u1	99
29) Benzo(g,h,i)perylene	27.592	276	14503	0.359	ng/u1	99

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

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