

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041323\
 Data File : BN024862.D
 Acq On : 13 Apr 2023 15:11
 Operator : CG/JU
 Sample : PB152092BS
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS092

Quant Time: Apr 14 02:20:30 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN040723.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 13 02:44:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.955	152	7643	0.400	ng/u1	0.00
4) Naphthalene-d8	10.763	136	22484	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.592	164	12325	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.338	188	27489	0.400	ng/u1	0.00
17) Chrysene-d12	21.531	240	22872	0.400	ng/u1	0.00
23) Perylene-d12	23.986	264	19577	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	5668	0.679	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.353	152	10205	0.354	ng/u1	0.00
18) Fluoranthene-d10	19.367	212	22442	0.395	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.352	88	14758	1.605	ng/u1	94
5) Naphthalene	10.813	128	19945	0.348	ng/u1	99
7) 2-Methylnaphthalene	12.424	142	12193	0.342	ng/u1	99
8) 1-Methylnaphthalene	12.644	142	13044	0.345	ng/u1	99
10) Acenaphthylene	14.314	152	15967	0.345	ng/u1	99
11) Acenaphthene	14.657	153	13603	0.343	ng/u1	99
12) Fluorene	15.642	166	15320	0.345	ng/u1	99
14) Pentachlorophenol	16.991	266	4096	0.699	ng/u1	99
15) Phenanthrene	17.380	178	26306	0.340	ng/u1	100
16) Anthracene	17.477	178	21767	0.338	ng/u1	98
19) Fluoranthene	19.399	202	29850	0.379	ng/u1	98
20) Pyrene	19.757	202	31164	0.386	ng/u1	98
21) Benzo(a)anthracene	21.513	228	26368	0.365	ng/u1	100
22) Chrysene	21.568	228	28968	0.370	ng/u1	100
24) Benzo(b)fluoranthene	23.232	252	28947	0.369	ng/u1	97
25) Benzo(k)fluoranthene	23.284	252	27133	0.354	ng/u1	95
26) Benzo(a)pyrene	23.872	252	22994	0.353	ng/u1	94
27) Indeno(1,2,3-cd)pyrene	26.539	276	25629	0.334	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.563	278	19706	0.326	ng/u1	97
29) Benzo(g,h,i)perylene	27.334	276	22408	0.345	ng/u1	100

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

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