

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032924\
 Data File : BN030529.D
 Acq On : 29 Mar 2024 12:36
 Operator : MA/JU
 Sample : PB159848BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS848

Quant Time: Mar 29 13:01:05 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 27 15:11:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.704	152	6389	0.400	ng/u1	-0.02
4) Naphthalene-d8	10.486	136	18883	0.400	ng/u1	-0.02
9) Acenaphthene-d10	14.346	164	10546	0.400	ng/u1	-0.02
13) Phenanthrene-d10	17.092	188	23598	0.400	ng/u1	-0.02
17) Chrysene-d12	21.287	240	21790	0.400	ng/u1	-0.01
23) Perylene-d12	23.532	264	21475	0.400	ng/u1	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.202	96	4947	0.707	ng/u1	-0.02
6) 2-Methylnaphthalene-d10	12.081	152	10388	0.365	ng/u1	-0.02
18) Fluoranthene-d10	19.124	212	22980	0.341	ng/u1	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.236	88	13616	1.919	ng/u1#	86
5) Naphthalene	10.536	128	18871	0.367	ng/u1	100
7) 2-Methylnaphthalene	12.158	142	12271	0.361	ng/u1	100
8) 1-Methylnaphthalene	12.378	142	12389	0.344	ng/u1	100
10) Acenaphthylene	14.063	152	17744	0.373	ng/u1	99
11) Acenaphthene	14.410	153	13519	0.365	ng/u1	97
12) Fluorene	15.400	166	15423	0.353	ng/u1	99
14) Pentachlorophenol	16.737	266	3507	0.813	ng/u1	100
15) Phenanthrene	17.134	178	26252	0.371	ng/u1	100
16) Anthracene	17.227	178	22407	0.376	ng/u1	99
19) Fluoranthene	19.157	202	30411	0.338	ng/u1	100
20) Pyrene	19.519	202	31284	0.354	ng/u1	99
21) Benzo(a)anthracene	21.273	228	29744	0.396	ng/u1	99
22) Chrysene	21.322	228	31699	0.382	ng/u1	100
24) Benzo(b)fluoranthene	22.854	252	31587	0.361	ng/u1	99
25) Benzo(k)fluoranthene	22.900	252	33777	0.366	ng/u1	99
26) Benzo(a)pyrene	23.433	252	28366	0.400	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	25.800	276	36751	0.458	ng/u1#	99
28) Dibenzo(a,h)anthracene	25.817	278	29381	0.461	ng/u1	98
29) Benzo(g,h,i)perylene	26.491	276	29897	0.439	ng/u1	98

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

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