

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071224\
 Data File : BN032653.D
 Acq On : 12 Jul 2024 01:34
 Operator : MA/JU
 Sample : PB161826BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS826

Quant Time: Jul 12 03:20:24 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN071224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 01:40:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.561	152	5535	0.400	ng/ul	-0.02
4) Naphthalene-d8	10.332	136	16618	0.400	ng/ul	-0.03
9) Acenaphthene-d10	14.207	164	8619	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.970	188	15334	0.400	ng/ul	-0.02
17) Chrysene-d12	21.179	240	13848	0.400	ng/ul	-0.01
23) Perylene-d12	23.374	264	16810	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.106	96	4436	0.667	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.933	152	8570	0.341	ng/ul	-0.04
18) Fluoranthene-d10	19.008	212	15976	0.354	ng/ul	-0.02
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.135	88	12190	1.610	ng/ul	91
5) Naphthalene	10.382	128	16367	0.367	ng/ul	99
7) 2-Methylnaphthalene	12.010	142	9744	0.350	ng/ul	98
8) 1-Methylnaphthalene	12.230	142	9681	0.336	ng/ul	98
10) Acenaphthylene	13.929	152	14937	0.429	ng/ul	99
11) Acenaphthene	14.272	153	10743	0.373	ng/ul	99
12) Fluorene	15.271	166	11097	0.337	ng/ul	99
14) Pentachlorophenol	16.640	266	2505	0.931	ng/ul	99
15) Phenanthrene	17.012	178	15792	0.378	ng/ul	100
16) Anthracene	17.109	178	14868	0.436	ng/ul	100
19) Fluoranthene	19.036	202	19926	0.354	ng/ul	99
20) Pyrene	19.403	202	20865	0.358	ng/ul	99
21) Benzo(a)anthracene	21.162	228	19197	0.436	ng/ul	96
22) Chrysene	21.214	228	22285	0.349	ng/ul	96
24) Benzo(b)fluoranthene	22.713	252	20646	0.348	ng/ul	90
25) Benzo(k)fluoranthene	22.757	252	24017	0.332	ng/ul#	91
26) Benzo(a)pyrene	23.278	252	21609	0.421	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.586	276	27395	0.394	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.596	278	21852	0.406	ng/ul#	94
29) Benzo(g,h,i)perylene	26.263	276	22489	0.387	ng/ul	96

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

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