

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081522\
 Data File : BN021230.D
 Acq On : 15 Aug 2022 11:07
 Operator : CG/JU
 Sample : PB146961BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS961

Quant Time: Aug 15 11:57:34 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN080322.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Aug 13 04:02:44 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.797	152	2882	0.400 ng/u1	0.00
4) Naphthalene-d8	10.584	136	10542	0.400 ng/u1	0.00
9) Acenaphthene-d10	14.437	164	5902	0.400 ng/u1	0.00
13) Phenanthrene-d10	17.191	188	8984	0.400 ng/u1	0.00
17) Chrysene-d12	21.395	240	5717	0.400 ng/u1	0.00
23) Perylene-d12	23.725	264	4602	0.400 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.194	96	2577	0.699 ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.190	152	5252	0.340 ng/u1	0.00
18) Fluoranthene-d10	19.229	212	8228	0.466 ng/u1	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.228	88	6460	1.701 ng/u1	87
5) Naphthalene	10.633	128	10193	0.315 ng/u1	99
7) 2-Methylnaphthalene	12.261	142	5974	0.304 ng/u1	99
8) 1-Methylnaphthalene	12.476	142	6583	0.335 ng/u1	100
10) Acenaphthylene	14.155	152	8366	0.303 ng/u1	99
11) Acenaphthene	14.497	153	7088	0.311 ng/u1	99
12) Fluorene	15.496	166	7486	0.300 ng/u1	100
14) Pentachlorophenol	16.862	266	514	0.371 ng/u1	99
15) Phenanthrene	17.229	178	9766	0.304 ng/u1	100
16) Anthracene	17.331	178	7356	0.295 ng/u1	99
19) Fluoranthene	19.257	202	10113	0.400 ng/u1	99
20) Pyrene	19.620	202	10495	0.402 ng/u1	95
21) Benzo(a)anthracene	21.381	228	5519	0.296 ng/u1	100
22) Chrysene	21.433	228	7614	0.310 ng/u1	100
24) Benzo(b)fluoranthene	23.017	252	6797	0.319 ng/u1	91
25) Benzo(k)fluoranthene	23.064	252	7279	0.324 ng/u1	96
26) Benzo(a)pyrene	23.625	252	6477	0.333 ng/u1#	73
27) Indeno(1,2,3-cd)pyrene	26.136	276	6857	0.295 ng/u1#	95
28) Dibenzo(a,h)anthracene	26.159	278	5079	0.277 ng/u1	94
29) Benzo(g,h,i)perylene	26.867	276	6363	0.307 ng/u1	97

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

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