

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100523\
 Data File : BN028108.D
 Acq On : 07 Oct 2023 06:16
 Operator : MA/JU
 Sample : PB156033BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS033

Quant Time: Oct 08 00:50:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN100323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 06 05:15:40 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.937	152	4682	0.400	ng/ul	0.00
4) Naphthalene-d8	10.751	136	14432	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.578	164	7173	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.337	188	13624	0.400	ng/ul	0.00
17) Chrysene-d12	21.521	240	12118	0.400	ng/ul	0.00
23) Perylene-d12	23.985	264	13439	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.326	96	4642	0.789	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.335	152	9061	0.390	ng/ul	0.00
18) Fluoranthene-d10	19.362	212	16357	0.426	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.364	88	13932	2.238	ng/ul#	86
5) Naphthalene	10.801	128	15798	0.386	ng/ul	99
7) 2-Methylnaphthalene	12.407	142	9494	0.367	ng/ul	98
8) 1-Methylnaphthalene	12.627	142	9285	0.347	ng/ul	100
10) Acenaphthylene	14.310	152	13643	0.419	ng/ul	99
11) Acenaphthene	14.643	153	10076	0.389	ng/ul	99
12) Fluorene	15.628	166	10731	0.372	ng/ul	98
14) Pentachlorophenol	16.974	266	2505	0.857	ng/ul	100
15) Phenanthrene	17.375	178	16110	0.397	ng/ul	99
16) Anthracene	17.472	178	14775	0.407	ng/ul	98
19) Fluoranthene	19.394	202	18266	0.395	ng/ul	99
20) Pyrene	19.761	202	19096	0.398	ng/ul	100
21) Benzo(a)anthracene	21.504	228	18704	0.423	ng/ul	100
22) Chrysene	21.559	228	18437	0.413	ng/ul	99
24) Benzo(b)fluoranthene	23.222	252	18915	0.380	ng/ul	94
25) Benzo(k)fluoranthene	23.272	252	19942	0.402	ng/ul	95
26) Benzo(a)pyrene	23.874	252	18317	0.431	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.575	276	23242	0.452	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.592	278	18825	0.450	ng/ul	97
29) Benzo(g,h,i)perylene	27.387	276	19053	0.453	ng/ul	98

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

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