

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110722\
 Data File : BN022553.D
 Acq On : 08 Nov 2022 08:11
 Operator : CG/JU
 Sample : PB148828BS
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS828

Quant Time: Nov 09 01:15:29 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 07 23:39:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.911	152	1613	0.400	ng/ul	0.00
4) Naphthalene-d8	10.739	136	5722	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.591	164	3828	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.349	188	8328	0.400	ng/ul	0.00
17) Chrysene-d12	21.553	240	7318	0.400	ng/ul	0.00
23) Perylene-d12	24.002	264	6063	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	1549	0.739	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.340	152	4037	0.466	ng/ul	0.00
18) Fluoranthene-d10	19.384	212	10088	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.257	88	3900	1.812	ng/ul#	89
5) Naphthalene	10.789	128	6819	0.403	ng/ul	98
7) 2-Methylnaphthalene	12.411	142	4570	0.429	ng/ul	100
8) 1-Methylnaphthalene	12.631	142	5012	0.457	ng/ul	100
10) Acenaphthylene	14.309	152	7245	0.422	ng/ul	98
11) Acenaphthene	14.651	153	5694	0.393	ng/ul	100
12) Fluorene	15.646	166	6455	0.383	ng/ul	99
14) Pentachlorophenol	17.007	266	1487	0.751	ng/ul	99
15) Phenanthrene	17.391	178	10767	0.385	ng/ul	99
16) Anthracene	17.484	178	9779	0.415	ng/ul	99
19) Fluoranthene	19.417	202	12546	0.374	ng/ul	97
20) Pyrene	19.779	202	12944	0.384	ng/ul	98
21) Benzo(a)anthracene	21.538	228	11760	0.429	ng/ul	99
22) Chrysene	21.591	228	11804	0.406	ng/ul	99
24) Benzo(b)fluoranthene	23.248	252	11430	0.375	ng/ul	96
25) Benzo(k)fluoranthene	23.298	252	11036	0.373	ng/ul	98
26) Benzo(a)pyrene	23.891	252	9716	0.400	ng/ul	94
27) Indeno(1,2,3-cd)pyrene	26.568	276	11486	0.381	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.588	278	9271	0.385	ng/ul	96
29) Benzo(g,h,i)perylene	27.356	276	9321	0.379	ng/ul	97

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

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