

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071023\
 Data File : BN026439.D
 Acq On : 10 Jul 2023 09:42
 Operator : MA/JU
 Sample : PB153949BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS949

Quant Time: Jul 11 04:17:27 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	3803	0.400	ng/ul	0.00
4) Naphthalene-d8	10.732	136	13449	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.566	164	7234	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.326	188	8852	0.400	ng/ul #	0.00
17) Chrysene-d12	21.523	240	6722	0.400	ng/ul	0.00
23) Perylene-d12	23.981	264	6553	0.400	ng/ul #	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	2891	0.675	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.332	152	5459	0.264	ng/ul	0.00
18) Fluoranthene-d10	19.354	212	8284	0.324	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.331	88	7576	1.666	ng/ul#	84
5) Naphthalene	10.781	128	12964	0.336	ng/ul	98
7) 2-Methylnaphthalene	12.404	142	6876	0.280	ng/ul	99
8) 1-Methylnaphthalene	12.613	142	7869	0.309	ng/ul	100
10) Acenaphthylene	14.293	152	11379	0.336	ng/ul	100
11) Acenaphthene	14.630	153	9441	0.341	ng/ul	100
12) Fluorene	15.625	166	9076	0.286	ng/ul	100
14) Pentachlorophenol	16.992	266	1411	0.445	ng/ul	98
15) Phenanthrene	17.368	178	9238	0.316	ng/ul	97
16) Anthracene	17.465	178	7414	0.285	ng/ul#	94
19) Fluoranthene	19.387	202	12533	0.349	ng/ul	98
20) Pyrene	19.749	202	13576	0.364	ng/ul	96
21) Benzo(a)anthracene	21.509	228	8492	0.306	ng/ul	99
22) Chrysene	21.564	228	9817	0.340	ng/ul	99
24) Benzo(b)fluoranthene	23.230	252	9876	0.353	ng/ul	90
25) Benzo(k)fluoranthene	23.283	252	10423	0.371	ng/ul	92
26) Benzo(a)pyrene	23.873	252	8482	0.349	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.558	276	10290	0.388	ng/ul#	89
28) Dibenzo(a,h)anthracene	26.571	278	7660	0.381	ng/ul#	92
29) Benzo(g,h,i)perylene	27.349	276	9680	0.410	ng/ul	97

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

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