

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN100423\
 Data File : BN028039.D
 Acq On : 04 Oct 2023 17:33
 Operator : MA/JU
 Sample : PB155913BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS913

Quant Time: Oct 05 02:05:33 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 : Tue Oct 03 15:50:04 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	4755	0.400	ng/ul	0.00
4) Naphthalene-d8	10.762	136	15838	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.587	164	9518	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.342	188	20289	0.400	ng/ul	0.00
17) Chrysene-d12	21.527	240	13627	0.400	ng/ul	0.00
23) Perylene-d12	23.997	264	13931	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.325	96	3935	0.659	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	9888	0.387	ng/ul	0.00
18) Fluoranthene-d10	19.366	212	20570	0.476	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.359	88	9598	1.518	ng/ul	90
5) Naphthalene	10.812	128	14872	0.331	ng/ul	100
7) 2-Methylnaphthalene	12.418	142	9731	0.343	ng/ul	100
8) 1-Methylnaphthalene	12.638	142	10262	0.349	ng/ul	100
10) Acenaphthylene	14.319	152	14918	0.346	ng/ul	100
11) Acenaphthene	14.652	153	11385	0.331	ng/ul	99
12) Fluorene	15.637	166	12857	0.336	ng/ul	98
14) Pentachlorophenol	16.978	266	2985	0.686	ng/ul	99
15) Phenanthrene	17.384	178	20188	0.334	ng/ul	100
16) Anthracene	17.477	178	18292	0.338	ng/ul	100
19) Fluoranthene	19.399	202	21441	0.412	ng/ul	99
20) Pyrene	19.766	202	22024	0.408	ng/ul	99
21) Benzo(a)anthracene	21.510	228	17380	0.349	ng/ul	99
22) Chrysene	21.565	228	17055	0.339	ng/ul	98
24) Benzo(b)fluoranthene	23.231	252	17259	0.334	ng/ul	90
25) Benzo(k)fluoranthene	23.281	252	17200	0.335	ng/ul#	90
26) Benzo(a)pyrene	23.886	252	14959	0.340	ng/ul#	89
27) Indeno(1,2,3-cd)pyrene	26.592	276	21028	0.395	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.609	278	17248	0.398	ng/ul	95
29) Benzo(g,h,i)perylene	27.407	276	17941	0.412	ng/ul#	93

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

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