

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025031.D
 Acq On : 20 Apr 2023 12:27
 Operator : CG/JU
 Sample : 02296-02
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCBJ7

Quant Time: Apr 21 04:17:28 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN042023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 21 04:13:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.925	152	6567	0.400	ng/u1	0.00
4) Naphthalene-d8	10.737	136	20069	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.575	164	10875	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.326	188	23552	0.400	ng/u1	0.00
17) Chrysene-d12	21.520	240	18552	0.400	ng/u1	0.00
23) Perylene-d12	23.970	264	15540	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.301	96	3806	0.560	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.326	152	7762	0.301	ng/u1	0.00
18) Fluoranthene-d10	19.354	212	17390	0.355	ng/u1	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.787	128	15322	0.298	ng/u1	99
7) 2-Methylnaphthalene	12.403	142	17365	0.551	ng/u1	100
10) Acenaphthylene	14.297	152	9764	0.241	ng/u1	100
12) Fluorene	15.625	166	19090	0.500	ng/u1	99
14) Pentachlorophenol	16.975	266	4025	0.870	ng/u1	100
15) Phenanthrene	17.364	178	20441	0.307	ng/u1	99
16) Anthracene	17.461	178	25549	0.470	ng/u1	100
19) Fluoranthene	19.387	202	19878	0.296	ng/u1	99
20) Pyrene	19.749	202	25829	0.375	ng/u1	99
21) Benzo(a)anthracene	21.505	228	26483	0.474	ng/u1	99
22) Chrysene	21.558	228	29749	0.488	ng/u1	100
24) Benzo(b)fluoranthene	23.218	252	46656	0.740	ng/u1	97
25) Benzo(k)fluoranthene	23.268	252	20260	0.334	ng/u1	97
26) Benzo(a)pyrene	23.861	252	18826	0.361	ng/u1#	93
28) Dibenzo(a,h)anthracene	26.541	278	37156	0.791	ng/u1	97
29) Benzo(g,h,i)perylene	27.305	276	17923	0.351	ng/u1	99

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

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