

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042023\
 Data File : BN025034.D
 Acq On : 20 Apr 2023 14:26
 Operator : CG/JU
 Sample : 02296-13MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCBL6MSD

Quant Time: Apr 21 04:19:47 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	7.930	152	8876	0.400	ng/u1	0.00
4) Naphthalene-d8	10.743	136	27904	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.575	164	15877	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.322	188	34645	0.400	ng/u1	0.00
17) Chrysene-d12	21.517	240	25510	0.400	ng/u1	0.00
23) Perylene-d12	23.964	264	20627	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	2261	0.246	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.327	152	12787	0.357	ng/u1	0.00
18) Fluoranthene-d10	19.350	212	32402	0.482	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.344	88	6477	0.654	ng/u1	88
5) Naphthalene	10.792	128	27215	0.381	ng/u1	100
7) 2-Methylnaphthalene	12.404	142	15838	0.362	ng/u1	100
8) 1-Methylnaphthalene	12.618	142	16419	0.352	ng/u1	100
10) Acenaphthylene	14.293	152	21682	0.367	ng/u1	99
11) Acenaphthene	14.635	153	18590	0.373	ng/u1	99
12) Fluorene	15.620	166	21248	0.381	ng/u1	99
14) Pentachlorophenol	16.975	266	5212	0.766	ng/u1	100
15) Phenanthrene	17.364	178	37702	0.386	ng/u1	100
16) Anthracene	17.457	178	30329	0.379	ng/u1	100
19) Fluoranthene	19.382	202	43998	0.477	ng/u1	100
20) Pyrene	19.745	202	45003	0.475	ng/u1	100
21) Benzo(a)anthracene	21.503	228	34204	0.445	ng/u1	100
22) Chrysene	21.555	228	37571	0.448	ng/u1	99
24) Benzo(b)fluoranthene	23.215	252	34112	0.408	ng/u1	98
25) Benzo(k)fluoranthene	23.262	252	33855	0.421	ng/u1	99
26) Benzo(a)pyrene	23.859	252	26293	0.380	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.514	276	32566	0.408	ng/u1#	98
28) Dibenzo(a,h)anthracene	26.537	278	25678	0.412	ng/u1	98
29) Benzo(g,h,i)perylene	27.292	276	28726	0.423	ng/u1	99

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

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