

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN033023\
 Data File : BN024657.D
 Acq On : 30 Mar 2023 23:50
 Operator : CG/JU
 Sample : 02038-10MSD
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 EYF73MSD

Quant Time: Mar 31 02:13:54 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN033023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 31 02:10:21 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.000	152	8038	0.400	ng/u1	0.00
4) Naphthalene-d8	10.812	136	23595	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.638	164	12651	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.384	188	27466	0.400	ng/u1	0.00
17) Chrysene-d12	21.568	240	21237	0.400	ng/u1	0.00
23) Perylene-d12	24.035	264	17707	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.376	96	2330	0.282	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.396	152	11609	0.399	ng/u1	0.00
18) Fluoranthene-d10	19.408	212	27852	0.536	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.410	88	6956	0.801	ng/u1	97
5) Naphthalene	10.861	128	24508	0.423	ng/u1	99
7) 2-Methylnaphthalene	12.473	142	15095	0.418	ng/u1	100
8) 1-Methylnaphthalene	12.687	142	15838	0.420	ng/u1	100
10) Acenaphthylene	14.361	152	18953	0.424	ng/u1	100
11) Acenaphthene	14.698	153	16837	0.434	ng/u1	98
12) Fluorene	15.684	166	19105	0.431	ng/u1	99
14) Pentachlorophenol	17.033	266	5325	0.897	ng/u1	98
15) Phenanthrene	17.426	178	33885	0.450	ng/u1	100
16) Anthracene	17.515	178	27449	0.436	ng/u1	100
19) Fluoranthene	19.436	202	40251	0.538	ng/u1	98
20) Pyrene	19.799	202	41200	0.541	ng/u1	97
21) Benzo(a)anthracene	21.554	228	34019	0.533	ng/u1	99
22) Chrysene	21.606	228	37680	0.537	ng/u1	99
24) Benzo(b)fluoranthene	23.281	252	37217	0.506	ng/u1	99
25) Benzo(k)fluoranthene	23.328	252	35917	0.507	ng/u1	99
26) Benzo(a)pyrene	23.924	252	29459	0.509	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	26.603	276	34764	0.504	ng/u1#	99
28) Dibenzo(a,h)anthracene	26.623	278	27458	0.498	ng/u1	98
29) Benzo(g,h,i)perylene	27.404	276	28860	0.487	ng/u1	98

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

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