

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN081523\
 Data File : BN026903.D
 Acq On : 15 Aug 2023 16:18
 Operator : MA/JU
 Sample : 03965-02MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4832MS

Quant Time: Aug 16 00:54:08 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN080323.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 16 00:50:05 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.110	152	12985	0.400	ng/ul	0.00
4) Naphthalene-d8	10.938	136	40534	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.735	164	25050	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.481	188	59121	0.400	ng/ul	0.00
17) Chrysene-d12	21.656	240	53123	0.400	ng/ul	0.00
23) Perylene-d12	24.208	264	53293	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.439	96	6462	0.419	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.511	152	18187	0.305	ng/ul	0.00
18) Fluoranthene-d10	19.497	212	52222	0.311	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.477	88	15457	0.982	ng/ul#	91
5) Naphthalene	10.988	128	30383	0.265	ng/ul	99
7) 2-Methylnaphthalene	12.583	142	19653	0.283	ng/ul	100
8) 1-Methylnaphthalene	12.797	142	21083	0.290	ng/ul	100
10) Acenaphthylene	14.467	152	31531	0.225	ng/ul	99
11) Acenaphthene	14.800	153	24585	0.253	ng/ul	99
12) Fluorene	15.781	166	28943	0.263	ng/ul	98
15) Phenanthrene	17.523	178	125068	0.644	ng/ul	100
16) Anthracene	17.616	178	47614	0.271	ng/ul	100
19) Fluoranthene	19.529	202	283808	1.286	ng/ul	100
20) Pyrene	19.892	202	260716	1.107	ng/ul	99
21) Benzo(a)anthracene	21.638	228	151793	0.718	ng/ul	100
22) Chrysene	21.697	228	155632	0.746	ng/ul	99
24) Benzo(b)fluoranthene	23.416	252	218622	1.085	ng/ul	99
25) Benzo(k)fluoranthene	23.465	252	106353	0.535	ng/ul	100
26) Benzo(a)pyrene	24.091	252	136273	0.749	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.898	276	132994	0.559	ng/ul#	95
28) Dibenzo(a,h)anthracene	26.931	278	60991	0.322	ng/ul	100
29) Benzo(g,h,i)perylene	27.739	276	103366	0.522	ng/ul	100

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

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