

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070723\
 Data File : BN026356.D
 Acq On : 07 Jul 2023 13:16
 Operator : MA/JU
 Sample : 03284-10MS
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCGZ6MS

Quant Time: Jul 07 15:53:11 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 07 14:46:07 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.938	152	4104	0.400	ng/ul	0.00
4) Naphthalene-d8	10.743	136	16250	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.575	164	9033	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.330	188	15528	0.400	ng/ul	0.00
17) Chrysene-d12	21.535	240	9208	0.400	ng/ul	0.00
23) Perylene-d12	23.981	264	9618	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.305	96	9885	2.139	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.337	152	7698	0.308	ng/ul	0.00
18) Fluoranthene-d10	19.359	212	14514	0.414	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.339	88	5819	1.186	ng/ul#	91
5) Naphthalene	10.792	128	16403	0.352	ng/ul	99
7) 2-Methylnaphthalene	12.409	142	9883	0.333	ng/ul	99
8) 1-Methylnaphthalene	12.623	142	10504	0.342	ng/ul	100
10) Acenaphthylene	14.297	152	14808	0.350	ng/ul	99
11) Acenaphthene	14.640	153	11677	0.337	ng/ul	98
12) Fluorene	15.630	166	12617	0.318	ng/ul	99
14) Pentachlorophenol	16.992	266	2370	0.426	ng/ul	99
15) Phenanthrene	17.372	178	19163	0.374	ng/ul	99
16) Anthracene	17.465	178	15993	0.351	ng/ul	98
19) Fluoranthene	19.387	202	19944	0.406	ng/ul	99
20) Pyrene	19.754	202	20874	0.408	ng/ul	97
21) Benzo(a)anthracene	21.517	228	14164	0.373	ng/ul	99
22) Chrysene	21.570	228	15358	0.389	ng/ul	98
24) Benzo(b)fluoranthene	23.236	252	14692	0.358	ng/ul	93
25) Benzo(k)fluoranthene	23.283	252	15596	0.378	ng/ul	95
26) Benzo(a)pyrene	23.876	252	12951	0.363	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.544	276	17061	0.438	ng/ul#	96
28) Dibenzo(a,h)anthracene	26.554	278	13215	0.448	ng/ul	98
29) Benzo(g,h,i)perylene	27.329	276	15396	0.444	ng/ul	97

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

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