

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071023\
 Data File : BN026441.D
 Acq On : 10 Jul 2023 11:04
 Operator : MA/JU
 Sample : 03404-06MS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW43MS

Quant Time: Jul 11 04:17:53 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 : Sun Jul 09 15:51:14 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.921	152	4069	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.730	136	14094	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.565	164	7263	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.321	188	11406	0.400	ng/ul	0.00
17) Chrysene-d12	21.519	240	7610	0.400	ng/ul	0.00
23) Perylene-d12	23.963	264	7450	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.297	96	2362	0.515	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.325	152	4925	0.227	ng/ul	0.00
18) Fluoranthene-d10	19.349	212	7693	0.266	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.331	88	5725	1.177	ng/ul#	89
5) Naphthalene	10.780	128	10612	0.263	ng/ul	97
7) 2-Methylnaphthalene	12.402	142	5780	0.225	ng/ul	99
8) 1-Methylnaphthalene	12.611	142	6347	0.238	ng/ul	99
10) Acenaphthylene	14.288	152	9058	0.266	ng/ul	98
11) Acenaphthene	14.625	153	7286	0.262	ng/ul	99
12) Fluorene	15.620	166	7211	0.226	ng/ul	98
14) Pentachlorophenol	16.983	266	652	0.159	ng/ul	95
15) Phenanthrene	17.363	178	9818	0.261	ng/ul	96
16) Anthracene	17.456	178	9528	0.285	ng/ul	95
19) Fluoranthene	19.381	202	10644	0.262	ng/ul	99
20) Pyrene	19.744	202	11475	0.271	ng/ul	99
21) Benzo(a)anthracene	21.505	228	7842	0.250	ng/ul	98
22) Chrysene	21.554	228	8810	0.270	ng/ul	99
24) Benzo(b)fluoranthene	23.220	252	8453	0.266	ng/ul	86
25) Benzo(k)fluoranthene	23.267	252	9367	0.293	ng/ul#	86
26) Benzo(a)pyrene	23.858	252	7222	0.261	ng/ul#	81
27) Indeno(1,2,3-cd)pyrene	26.523	276	8977	0.298	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.540	278	6890	0.302	ng/ul#	88
29) Benzo(g,h,i)perylene	27.308	276	8266	0.308	ng/ul	96

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

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