

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070723\
 Data File : BN026371.D
 Acq On : 07 Jul 2023 23:11
 Operator : MA/JU
 Sample : 03237-05MSD
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YBW77MSD

Quant Time: Jul 08 04:47:23 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	4567	0.400	ng/ul	0.00
4) Naphthalene-d8	10.748	136	16349	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.575	164	9832	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.330	188	14399	0.400	ng/ul	0.00
17) Chrysene-d12	21.523	240	9221	0.400	ng/ul	0.00
23) Perylene-d12	23.967	264	9919	0.400	ng/ul	#-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.310	96	796	0.155	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.337	152	5796	0.230	ng/ul	0.00
18) Fluoranthene-d10	19.359	212	10836	0.309	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.343	88	1973	0.361	ng/ul	95
5) Naphthalene	10.798	128	11895	0.254	ng/ul	99
7) 2-Methylnaphthalene	12.414	142	6480	0.217	ng/ul	98
8) 1-Methylnaphthalene	12.623	142	7277	0.235	ng/ul	99
10) Acenaphthylene	14.302	152	11215	0.244	ng/ul	99
11) Acenaphthene	14.640	153	9976	0.265	ng/ul	96
12) Fluorene	15.630	166	8214	0.190	ng/ul	99
14) Pentachlorophenol	16.996	266	1559	0.302	ng/ul	98
15) Phenanthrene	17.372	178	13530	0.284	ng/ul	97
16) Anthracene	17.469	178	11178	0.264	ng/ul	96
19) Fluoranthene	19.387	202	15139	0.307	ng/ul	100
20) Pyrene	19.754	202	15886	0.310	ng/ul	96
21) Benzo(a)anthracene	21.506	228	11132	0.293	ng/ul	98
22) Chrysene	21.561	228	11848	0.299	ng/ul	98
24) Benzo(b)fluoranthene	23.218	252	12470	0.295	ng/ul	88
25) Benzo(k)fluoranthene	23.271	252	12752	0.300	ng/ul#	90
26) Benzo(a)pyrene	23.862	252	10962	0.298	ng/ul#	84
27) Indeno(1,2,3-cd)pyrene	26.524	276	14598	0.363	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.541	278	11169	0.367	ng/ul	94
29) Benzo(g,h,i)perylene	27.312	276	13088	0.366	ng/ul	98

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

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