

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072023\
 Data File : BN026617.D
 Acq On : 21 Jul 2023 07:10
 Operator : MA/JU
 Sample : 03472-30MS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A46Q2MS

Quant Time: Jul 21 07:40:17 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN071223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jul 20 22:32:26 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.836	152	1705	0.400	ng/ul	-0.03
4) Naphthalene-d8	10.636	136	6021	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.472	164	3803	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.219	188	9610	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.411	240	9794	0.400	ng/ul	0.00
23) Perylene-d12	23.784	264	9810	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.203	96	771	0.343	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.231	152	2579	0.315	ng/ul	-0.02
18) Fluoranthene-d10	19.246	212	10350	0.305	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.237	88	2041	0.961	ng/ul#	82
5) Naphthalene	10.691	128	5048	0.305	ng/ul	98
7) 2-Methylnaphthalene	12.302	142	3433	0.366	ng/ul	98
8) 1-Methylnaphthalene	12.517	142	3845	0.383	ng/ul	99
10) Acenaphthylene	14.194	152	6487	0.348	ng/ul	99
11) Acenaphthene	14.532	153	4778	0.337	ng/ul	99
12) Fluorene	15.526	166	6246	0.420	ng/ul	99
14) Pentachlorophenol	16.886	266	940	0.388	ng/ul	99
15) Phenanthrene	17.261	178	18977	0.620	ng/ul	98
16) Anthracene	17.358	178	6417	0.242	ng/ul	98
19) Fluoranthene	19.278	202	34658	0.755	ng/ul	98
20) Pyrene	19.641	202	42496	0.852	ng/ul#	97
21) Benzo(a)anthracene	21.396	228	25399	0.706	ng/ul	98
22) Chrysene	21.449	228	30751	0.751	ng/ul	99
24) Benzo(b)fluoranthene	23.062	252	33843	0.886	ng/ul	90
25) Benzo(k)fluoranthene	23.109	252	23823	0.595	ng/ul#	89
26) Benzo(a)pyrene	23.682	252	8363	0.246	ng/ul#	91
27) Indeno(1,2,3-cd)pyrene	26.241	276	25556	0.573	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.247	278	16237	0.474	ng/ul#	90
29) Benzo(g,h,i)perylene	26.989	276	22289	0.566	ng/ul	96

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

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