

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070324\
 Data File : BN032436.D
 Acq On : 03 Jul 2024 14:32
 Operator : MA/JU
 Sample : P2964-01DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN9DL

Quant Time: Jul 03 14:57:07 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.603	152	5146	0.400	ng/ul	0.00
4) Naphthalene-d8	10.376	136	16689	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.248	164	10546	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.007	188	20446	0.400	ng/ul	0.00
17) Chrysene-d12	21.214	240	14936	0.400	ng/ul	# 0.00
23) Perylene-d12	23.433	264	17057	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	293	0.051	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.988	152	561	0.022	ng/ul	0.01
18) Fluoranthene-d10	19.045	212	1557	0.035	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.426	128	39113	0.886	ng/ul	99
10) Acenaphthylene	13.971	152	17276	0.394	ng/ul	99
11) Acenaphthene	14.313	153	17591	0.517	ng/ul	99
12) Fluorene	15.312	166	5973	0.147	ng/ul	99
15) Phenanthrene	17.050	178	70448	1.257	ng/ul	99
16) Anthracene	17.143	178	53778	1.076	ng/ul	100
19) Fluoranthene	19.078	202	14992	0.264	ng/ul#	94
20) Pyrene	19.440	202	22990	0.387	ng/ul	99
21) Benzo(a)anthracene	21.200	228	8210	0.152	ng/ul	94
22) Chrysene	21.252	228	46108	0.779	ng/ul	99
24) Benzo(b)fluoranthene	22.769	252	22996	0.310	ng/ul	85
25) Benzo(k)fluoranthene	22.813	252	5846	0.081	ng/ul#	39
26) Benzo(a)pyrene	23.339	252	28387	0.551	ng/ul#	94
27) Indeno(1,2,3-cd)pyrene	25.673	276	18623	0.231	ng/ul#	97
28) Dibenzo(a,h)anthracene	25.683	278	16307	0.272	ng/ul#	83
29) Benzo(g,h,i)perylene	26.360	276	45990	0.715	ng/ul	99

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

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