

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070524\
 Data File : BN032509.D
 Acq On : 05 Jul 2024 14:36
 Operator : MA/JU
 Sample : P2982-04
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DP5

Quant Time: Jul 05 15:02:53 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	7.595	152	4677	0.400	ng/ul	0.00
4) Naphthalene-d8	10.369	136	13838	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.238	164	7337	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.998	188	14098	0.400	ng/ul	0.00
17) Chrysene-d12	21.207	240	13067	0.400	ng/ul	# 0.00
23) Perylene-d12	23.428	264	15432	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	2271	0.437	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.970	152	5586	0.269	ng/ul	0.00
18) Fluoranthene-d10	19.035	212	12765	0.325	ng/ul	-0.01
Target Compounds					Qvalue	
5) Naphthalene	10.413	128	294522	8.042	ng/ul	99
10) Acenaphthylene	13.960	152	309127	10.130	ng/ul	99
11) Acenaphthene	14.302	153	389285	16.437	ng/ul	98
12) Fluorene	15.297	166	227627	8.066	ng/ul	99
15) Phenanthrene	17.040	178	604285	15.634	ng/ul	99
16) Anthracene	17.133	178	554159	16.074	ng/ul	99
19) Fluoranthene	19.067	202	432642	8.695	ng/ul	98
20) Pyrene	19.434	202	196732	3.783	ng/ul	99
21) Benzo(a)anthracene	21.192	228	61249	1.299	ng/ul	99
22) Chrysene	21.245	228	181937	3.515	ng/ul	99
24) Benzo(b)fluoranthene	22.759	252	195561	2.917	ng/ul	94
25) Benzo(k)fluoranthene	22.803	252	160953	2.471	ng/ul	96
26) Benzo(a)pyrene	23.329	252	267277	5.730	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	25.661	276	341619	4.677	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.665	278	324986	5.988	ng/ul	96
29) Benzo(g,h,i)perylene	26.349	276	246898	4.244	ng/ul	97

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

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