

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070124\
 Data File : BN032362.D
 Acq On : 01 Jul 2024 14:37
 Operator : MA/JU
 Sample : P2934-25DL2 20X
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN5DL2

Quant Time: Jul 01 15:02:14 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 : Thu Jun 27 13:23:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.612	152	6605	0.400	ng/ul	0.00
4) Naphthalene-d8	10.387	136	20639	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.258	164	11381	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.016	188	19910	0.400	ng/ul	0.00
17) Chrysene-d12	21.217	240	18001	0.400	ng/ul	0.00
23) Perylene-d12	23.436	264	21853	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	247	0.034	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.999	152	443	0.014	ng/ul	0.00
18) Fluoranthene-d10	19.050	212	863	0.016	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.437	128	49603	0.908	ng/ul	99
10) Acenaphthylene	13.976	152	40462	0.855	ng/ul	99
11) Acenaphthene	14.318	153	27399	0.746	ng/ul	98
12) Fluorene	15.317	166	14830	0.339	ng/ul	100
15) Phenanthrene	17.058	178	21276	0.390	ng/ul	99
16) Anthracene	17.151	178	13932	0.286	ng/ul	97
19) Fluoranthene	19.083	202	27686	0.404	ng/ul	99
20) Pyrene	19.445	202	34437	0.481	ng/ul	100
21) Benzo(a)anthracene	21.203	228	8970	0.138	ng/ul	96
22) Chrysene	21.255	228	17973	0.252	ng/ul	97
24) Benzo(b)fluoranthene	22.769	252	21600	0.228	ng/ul	88
25) Benzo(k)fluoranthene	22.813	252	18946	0.205	ng/ul#	84
26) Benzo(a)pyrene	23.342	252	24158	0.366	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.676	276	30490	0.295	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.686	278	19429	0.253	ng/ul#	90
29) Benzo(g,h,i)perylene	26.357	276	21220	0.258	ng/ul	94

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

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