

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070124\
 Data File : BN032360.D
 Acq On : 01 Jul 2024 13:24
 Operator : MA/JU
 Sample : P2934-25DL 10X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN5DL

Quant Time: Jul 01 13:49:31 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	6898	0.400	ng/ul	0.00
4) Naphthalene-d8	10.387	136	23345	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.258	164	13498	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.012	188	22526	0.400	ng/ul	-0.01
17) Chrysene-d12	21.220	240	16006	0.400	ng/ul	0.00
23) Perylene-d12	23.438	264	19146	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.152	96	457	0.060	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	1047	0.030	ng/ul	0.00
18) Fluoranthene-d10	19.050	212	1830	0.038	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.437	128	114792	1.858	ng/ul	99
10) Acenaphthylene	13.975	152	102681	1.829	ng/ul	99
11) Acenaphthene	14.318	153	66898	1.535	ng/ul	98
12) Fluorene	15.312	166	38179	0.735	ng/ul	99
15) Phenanthrene	17.054	178	51596	0.835	ng/ul	100
16) Anthracene	17.147	178	34680	0.630	ng/ul	100
19) Fluoranthene	19.078	202	60200	0.988	ng/ul	98
20) Pyrene	19.445	202	73040	1.147	ng/ul	98
21) Benzo(a)anthracene	21.205	228	18564	0.321	ng/ul	98
22) Chrysene	21.255	228	35408	0.558	ng/ul	98
24) Benzo(b)fluoranthene	22.769	252	42113	0.506	ng/ul	96
25) Benzo(k)fluoranthene	22.813	252	37528	0.464	ng/ul#	94
26) Benzo(a)pyrene	23.342	252	47047	0.813	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.673	276	59848	0.660	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.686	278	38427	0.571	ng/ul	97
29) Benzo(g,h,i)perylene	26.360	276	41279	0.572	ng/ul	98

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

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