

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN120225\
 Data File : BN038255.D
 Acq On : 02 Dec 2025 15:13
 Operator : RC/JU
 Sample : Q3688-03MSD
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 M-31MSD

Quant Time: Dec 03 00:36:10 2025
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN112625.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 01 16:36:13 2025
 Response via : Initial Calibration

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

Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.762	152	7837	0.400 ng/u1	0.00
4) Naphthalene-d8	10.567	136	23944	0.400 ng/u1	0.00
9) Acenaphthene-d10	14.427	164	11905	0.400 ng/u1	0.00
13) Phenanthrene-d10	17.179	188	17150	0.400 ng/u1	-0.01
17) Chrysene-d12	21.371	240	11866	0.400 ng/u1	0.00
23) Perylene-d12	23.678	264	12141	0.400 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.218	96	3284	0.399 ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.167	152	6696	0.196 ng/u1	0.00
18) Fluoranthene-d10	19.215	212	15450	0.335 ng/u1	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.248	88	8524	0.962 ng/u1	94
5) Naphthalene	10.616	128	20665	0.318 ng/u1	99
7) 2-Methylnaphthalene	12.238	142	13523	0.311 ng/u1	100
8) 1-Methylnaphthalene	12.458	142	13104	0.284 ng/u1	100
10) Acenaphthylene	14.144	152	19399	0.347 ng/u1	99
11) Acenaphthene	14.491	153	13927	0.335 ng/u1	100
12) Fluorene	15.481	166	15067	0.338 ng/u1	99
14) Pentachlorophenol	16.833	266	3691	1.342 ng/u1	97
15) Phenanthrene	17.221	178	22587	0.420 ng/u1	100
16) Anthracene	17.314	178	19555	0.429 ng/u1	100
19) Fluoranthene	19.243	202	23593	0.372 ng/u1	98
20) Pyrene	19.605	202	24199	0.385 ng/u1	97
21) Benzo(a)anthracene	21.354	228	19837	0.500 ng/u1	99
22) Chrysene	21.407	228	21957	0.446 ng/u1	99
24) Benzo(b)fluoranthene	22.979	252	21635	0.455 ng/u1	98
25) Benzo(k)fluoranthene	23.026	252	22929	0.458 ng/u1	96
26) Benzo(a)pyrene	23.575	252	19996	0.466 ng/u1	94
27) Indeno(1,2,3-cd)pyrene	26.028	276	26870	0.526 ng/u1#	98
28) Dibenzo(a,h)anthracene	26.045	278	20630	0.575 ng/u1	95
29) Benzo(g,h,i)perylene	26.746	276	22595	0.470 ng/u1	100

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

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