

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032524\
 Data File : BN030479.D
 Acq On : 25 Mar 2024 13:51
 Operator : MA/JU
 Sample : SSTDICV0.4
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SICV208

Quant Time: Mar 25 14:23:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN032524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 25 13:38:40 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.724	152	9948	0.400 ng/u1	0.00
4) Naphthalene-d8	10.512	136	28015	0.400 ng/u1	0.00
9) Acenaphthene-d10	14.362	164	15011	0.400 ng/u1	0.00
13) Phenanthrene-d10	17.111	188	33528	0.400 ng/u1	0.00
17) Chrysene-d12	21.305	240	24779	0.400 ng/u1	0.00
23) Perylene-d12	23.558	264	23779	0.400 ng/u1	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.218	96	4462	0.382 ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.101	152	15701	0.398 ng/u1	0.00
18) Fluoranthene-d10	19.146	212	31255	0.404 ng/u1	0.00
Target Compounds					
					Qvalue
2) 1,4-Dioxane	3.252	88	4955	0.410 ng/u1	97
5) Naphthalene	10.556	128	34241	0.429 ng/u1	100
7) 2-Methylnaphthalene	12.178	142	22621	0.437 ng/u1	100
8) 1-Methylnaphthalene	12.398	142	21155	0.386 ng/u1	100
10) Acenaphthylene	14.084	152	32345	0.425 ng/u1	99
11) Acenaphthene	14.427	153	24210	0.413 ng/u1	98
12) Fluorene	15.417	166	28003	0.417 ng/u1	98
14) Pentachlorophenol	16.752	266	2366	0.375 ng/u1	98
15) Phenanthrene	17.153	178	46522	0.413 ng/u1	100
16) Anthracene	17.242	178	38640	0.406 ng/u1	100
19) Fluoranthene	19.174	202	50148	0.425 ng/u1	98
20) Pyrene	19.536	202	50077	0.430 ng/u1	99
21) Benzo(a)anthracene	21.290	228	40424	0.396 ng/u1	99
22) Chrysene	21.343	228	45623	0.422 ng/u1	100
24) Benzo(b)fluoranthene	22.880	252	42862	0.381 ng/u1	99
25) Benzo(k)fluoranthene	22.927	252	47220	0.411 ng/u1	99
26) Benzo(a)pyrene	23.459	252	34196	0.380 ng/u1	99
27) Indeno(1,2,3-cd)pyrene	25.844	276	45078	0.413 ng/u1#	99
28) Dibenzo(a,h)anthracene	25.864	278	36580	0.419 ng/u1	97
29) Benzo(g,h,i)perylene	26.538	276	35687	0.378 ng/u1	99

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

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