

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013123\
 Data File : BN023869.D
 Acq On : 31 Jan 2023 19:53
 Operator : CG/JU
 Sample : PB150470BS
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS470

Quant Time: Feb 01 01:04:43 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN013023.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 01 01:00:14 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc Units	Dev(Min)
Internal Standards					
1) 1,4-Dichlorobenzene-d4	7.863	152	3578	0.400 ng/ul	0.00
4) Naphthalene-d8	10.666	136	9746	0.400 ng/ul	0.00
9) Acenaphthene-d10	14.512	164	6229	0.400 ng/ul	0.00
13) Phenanthrene-d10	17.256	188	14259	0.400 ng/ul	0.00
17) Chrysene-d12	21.450	240	13810	0.400 ng/ul	0.00
23) Perylene-d12	23.853	264	11934	0.400 ng/ul	0.00
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.247	96	2264	0.590 ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.261	152	5001	0.328 ng/ul	0.00
18) Fluoranthene-d10	19.286	212	13200	0.356 ng/ul	0.00
Target Compounds					
				Qvalue	
2) 1,4-Dioxane	3.281	88	6052	1.489 ng/ul#	81
5) Naphthalene	10.716	128	8425	0.316 ng/ul	99
7) 2-Methylnaphthalene	12.333	142	5535	0.319 ng/ul	100
8) 1-Methylnaphthalene	12.553	142	5900	0.332 ng/ul	99
10) Acenaphthylene	14.229	152	7372	0.309 ng/ul	100
11) Acenaphthene	14.572	153	6541	0.312 ng/ul	99
12) Fluorene	15.557	166	7780	0.307 ng/ul	97
14) Pentachlorophenol	16.909	266	2503	0.532 ng/ul	100
15) Phenanthrene	17.298	178	13386	0.315 ng/ul	100
16) Anthracene	17.391	178	11446	0.322 ng/ul	99
19) Fluoranthene	19.314	202	16531	0.360 ng/ul	98
20) Pyrene	19.681	202	16906	0.356 ng/ul	100
21) Benzo(a)anthracene	21.432	228	16426	0.350 ng/ul	99
22) Chrysene	21.485	228	17189	0.348 ng/ul	99
24) Benzo(b)fluoranthene	23.119	252	18333	0.348 ng/ul	99
25) Benzo(k)fluoranthene	23.168	252	17466	0.339 ng/ul	98
26) Benzo(a)pyrene	23.739	252	14683	0.334 ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.329	276	20164	0.327 ng/ul	100
28) Dibenzo(a,h)anthracene	26.346	278	16392	0.333 ng/ul	96
29) Benzo(g,h,i)perylene	27.091	276	17089	0.337 ng/ul	99

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

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