

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN013123\
 Data File : BN023880.D
 Acq On : 01 Feb 2023 02:38
 Operator : CG/JU
 Sample : PB150470BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS470

Quant Time: Feb 01 03:30:08 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	7.859	152	4622	0.400	ng/ul	0.00
4) Naphthalene-d8	10.667	136	12914	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.511	164	8300	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.255	188	19086	0.400	ng/ul	0.00
17) Chrysene-d12	21.446	240	18110	0.400	ng/ul	0.00
23) Perylene-d12	23.829	264	15687	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.248	96	2511	0.507	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	6033	0.299	ng/ul	0.00
18) Fluoranthene-d10	19.285	212	16308	0.335	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.281	88	6580	1.253	ng/ul#	81
5) Naphthalene	10.716	128	9942	0.281	ng/ul	99
7) 2-Methylnaphthalene	12.333	142	6575	0.286	ng/ul	99
8) 1-Methylnaphthalene	12.553	142	7073	0.301	ng/ul	100
10) Acenaphthylene	14.229	152	9307	0.293	ng/ul	99
11) Acenaphthene	14.571	153	7847	0.281	ng/ul	98
12) Fluorene	15.557	166	9321	0.276	ng/ul	97
14) Pentachlorophenol	16.909	266	3242	0.515	ng/ul	99
15) Phenanthrene	17.297	178	15772	0.277	ng/ul	99
16) Anthracene	17.390	178	13839	0.291	ng/ul	99
19) Fluoranthene	19.318	202	19200	0.319	ng/ul	99
20) Pyrene	19.676	202	19732	0.316	ng/ul	97
21) Benzo(a)anthracene	21.429	228	18996	0.309	ng/ul	99
22) Chrysene	21.481	228	19051	0.294	ng/ul	99
24) Benzo(b)fluoranthene	23.104	252	19248	0.278	ng/ul	97
25) Benzo(k)fluoranthene	23.150	252	18989	0.280	ng/ul	96
26) Benzo(a)pyrene	23.723	252	16255	0.282	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.295	276	19919	0.246	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.312	278	15962	0.246	ng/ul	96
29) Benzo(g,h,i)perylene	27.057	276	16475	0.247	ng/ul	98

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

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