

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032023\
 Data File : BN024405.D
 Acq On : 20 Mar 2023 11:08
 Operator : CG/JU
 Sample : PB151451BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS451

Manual Integrations
 APPROVED

Reviewed By : Christian Giraldo 03/21/2023
 Supervised By : Jagrut Upadhyay 03/21/2023

Quant Time: Mar 21 02:14:23 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN030923.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 21 02:12:19 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	8679	0.400	ng/u1	0.00
4) Naphthalene-d8	10.850	136	26909	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.670	164	12798	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.413	188	25290	0.400	ng/u1	0.00
17) Chrysene-d12	21.600	240	15954	0.400	ng/u1	# 0.00
23) Perylene-d12	24.103	264	11114m	0.400	ng/u1	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.397	96	5789	0.606	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.434	152	9457	0.296	ng/u1	0.00
18) Fluoranthene-d10	19.436	212	15014	0.354	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.435	88	14957	1.440	ng/u1	90
5) Naphthalene	10.900	128	21257	0.316	ng/u1	100
7) 2-Methylnaphthalene	12.506	142	12249	0.301	ng/u1	98
8) 1-Methylnaphthalene	12.726	142	12780	0.298	ng/u1	99
10) Acenaphthylene	14.393	152	16487	0.309	ng/u1	99
11) Acenaphthene	14.731	153	12887	0.313	ng/u1	98
12) Fluorene	15.716	166	13907	0.303	ng/u1	97
14) Pentachlorophenol	17.059	266	3138	0.447	ng/u1#	73
15) Phenanthrene	17.456	178	22412	0.309	ng/u1	100
16) Anthracene	17.548	178	18487	0.291	ng/u1	99
19) Fluoranthene	19.469	202	22664	0.357	ng/u1	96
20) Pyrene	19.831	202	22320	0.346	ng/u1	95
21) Benzo(a)anthracene	21.586	228	15505	0.298	ng/u1	100
22) Chrysene	21.641	228	17674	0.321	ng/u1	100
24) Benzo(b)fluoranthene	23.334	252	14796	0.349	ng/u1	89
25) Benzo(k)fluoranthene	23.392	252	14913	0.339	ng/u1#	89
26) Benzo(a)pyrene	23.989	252	11601	0.318	ng/u1#	84
27) Indeno(1,2,3-cd)pyrene	26.723	276	12230	0.297	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.750	278	8923	0.285	ng/u1	93
29) Benzo(g,h,i)perylene	27.532	276	11440	0.321	ng/u1	95

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

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