

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN032023\
 Data File : BN024407.D
 Acq On : 20 Mar 2023 12:20
 Operator : CG/JU
 Sample : PB151452BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS452

Quant Time: Mar 21 02:14:40 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	10081	0.400	ng/u1	0.00
4) Naphthalene-d8	10.850	136	31105	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.670	164	14765	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.413	188	28720	0.400	ng/u1	# 0.00
17) Chrysene-d12	21.600	240	16960	0.400	ng/u1	0.00
23) Perylene-d12	24.082	264	11249	0.400	ng/u1	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.401	96	6582	0.593	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.434	152	11050	0.299	ng/u1	0.00
18) Fluoranthene-d10	19.436	212	16655	0.369	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.435	88	17259	1.430	ng/u1	91
5) Naphthalene	10.900	128	24557	0.316	ng/u1	100
7) 2-Methylnaphthalene	12.506	142	14235	0.302	ng/u1	100
8) 1-Methylnaphthalene	12.726	142	14808	0.299	ng/u1	99
10) Acenaphthylene	14.393	152	18968	0.308	ng/u1	100
11) Acenaphthene	14.731	153	14964	0.315	ng/u1	98
12) Fluorene	15.716	166	15913	0.300	ng/u1	97
14) Pentachlorophenol	17.059	266	3426	0.429	ng/u1#	73
15) Phenanthrene	17.455	178	25398	0.309	ng/u1	99
16) Anthracene	17.548	178	20934	0.290	ng/u1	99
19) Fluoranthene	19.469	202	25140	0.372	ng/u1	97
20) Pyrene	19.831	202	24755	0.361	ng/u1	95
21) Benzo(a)anthracene	21.583	228	16319	0.295	ng/u1	100
22) Chrysene	21.635	228	18926	0.324	ng/u1	100
24) Benzo(b)fluoranthene	23.322	252	15652	0.364	ng/u1	92
25) Benzo(k)fluoranthene	23.372	252	15765	0.354	ng/u1#	90
26) Benzo(a)pyrene	23.971	252	12150	0.329	ng/u1#	84
27) Indeno(1,2,3-cd)pyrene	26.680	276	12877	0.309	ng/u1#	100
28) Dibenzo(a,h)anthracene	26.696	278	9402	0.297	ng/u1	96
29) Benzo(g,h,i)perylene	27.481	276	11861	0.328	ng/u1	96

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

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