

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN062422\
 Data File : BN020460.D
 Acq On : 26 Jun 2022 06:46
 Operator : CG/JU
 Sample : PB145701BS
 Misc :
 ALS Vial : 78 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS701

Quant Time: Jun 27 03:44:38 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062422.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 00:17:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.750	152	2436	0.400	ng/u1	0.00
4) Naphthalene-d8	10.529	136	8375	0.400	ng/u1	0.00
9) Acenaphthene-d10	14.377	164	5548	0.400	ng/u1	0.00
13) Phenanthrene-d10	17.120	188	12158	0.400	ng/u1	0.00
17) Chrysene-d12	21.308	240	10846	0.400	ng/u1	0.00
23) Perylene-d12	23.588	264	9078	0.400	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.201	96	2138	0.758	ng/u1	0.00
6) 2-Methylnaphthalene-d10	12.124	152	5511	0.401	ng/u1	0.00
18) Fluoranthene-d10	19.146	212	13431	0.376	ng/u1	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.235	88	5296	1.838	ng/u1#	80
5) Naphthalene	10.579	128	9341	0.361	ng/u1	99
7) 2-Methylnaphthalene	12.196	142	6217	0.359	ng/u1	99
8) 1-Methylnaphthalene	12.416	142	6612	0.403	ng/u1	99
10) Acenaphthylene	14.095	152	9437	0.359	ng/u1	99
11) Acenaphthene	14.437	153	7438	0.353	ng/u1	99
12) Fluorene	15.427	166	8709	0.349	ng/u1	100
14) Pentachlorophenol	16.786	266	576	0.235	ng/u1	98
15) Phenanthrene	17.158	178	14764	0.349	ng/u1	100
16) Anthracene	17.251	178	13027	0.357	ng/u1	100
19) Fluoranthene	19.179	202	16842	0.340	ng/u1	99
20) Pyrene	19.541	202	17515	0.341	ng/u1	100
21) Benzo(a)anthracene	21.291	228	14345	0.368	ng/u1	100
22) Chrysene	21.343	228	16049	0.370	ng/u1	100
24) Benzo(b)fluoranthene	22.898	252	15193	0.362	ng/u1	99
25) Benzo(k)fluoranthene	22.945	252	14852	0.350	ng/u1	99
26) Benzo(a)pyrene	23.488	252	14014	0.367	ng/u1	99
27) Indeno(1,2,3-cd)pyrene	25.921	276	14350	0.378	ng/u1#	98
28) Dibenzo(a,h)anthracene	25.938	278	11628	0.379	ng/u1	100
29) Benzo(g,h,i)perylene	26.639	276	12372	0.379	ng/u1	99

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

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